

Page 1

Received: 12/09/86

SPECTRIX S.O.G

REPORT

WORK ORDER # 86-12-V41

01/21/87 15:58:09

REPORT Keystone Env. Resources, Inc.
TO Water Quality Engineering
440 College Park Drive
Monroeville, PA 15146-1536
ATTEN Shannon Craig

PREPARED Spectrix Corporation

BY 3913 Fondren

Suite 100

Houston, Texas 77063-5821

CERTIFIED BY

ATTEN Sample Control

PHONE (713) 266-6800

CONTACT CRUZ

CLIENT KEYSTONE SAMPLES 7
COMPANY Keystone Company, Inc.
FACILITY _____

WORK ID Water
TAKEN Client
TRANS Client
TYPE Project 1573-73
P. O. # 14-5-50106
INVOICE under separate cover

Please call the above number if you have any questions.
NOTE: ALL SAMPLES WILL BE RETAINED FOR 90 DAYS AND THEN
DISCARDED. IF YOU WISH YOUR SAMPLES RETURNED TO YOUR FACILITY
CALL SAMPLE CONTROL AT THE ABOVE NUMBER.

SAMPLE IDENTIFICATION

- 01 LAB BLANK (LLW)
- 02 CAV-QW10-001 (LLW)
- 03 CAV-QW11-001 (LLW)
- 04 SCK-MW17-001 (LLW)
- 05 SCK-PD3-001 (LLW)
- 06 TRIP BLANK (LLW)
- 07 SCK-PD2-001 (LLW)

TEST CODES and NAMES used on this report

AG	Silver
AS	Argenic
BE	Berulium
CD	Cadmium
CN SU	Cyanide in Water
CR	Chromium
CU	Copper
FE	Iron
HG	Mercury
NI	Nickel
NO3 SU	Nitrate as N
PB	Lead
PEST W	EPA Pesticides/PCB's-Water
SB	Antimony
SE	Selenium
TL	Thallium
ZN	Zinc

003339

LABORATORY: SPECTRIX-HOUSTON
LABORATORY NO: 86-12-041
CLIENT: SPECTRIX-MONROEVILLE
SITE: SOUTH CAVALCADE

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 - Inorganic Matrix Duplicate and Matrix Spike

CHECKLIST

SECTION I - Support Information and QC Summaries

- Narrative
- Introduction to the Spectrix Organic Data Package
- Introduction to the Spectrix Inorganic Data Package
- Shipping Document, Chain-of-Custody
- ~~Organic MS/MSD Data Summary Form~~

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SPECTRIX
NARRATIVE - ORGANIC ANALYSES

CLIENT: SPECTRIX - MONROEVILLE
LAB NO: 8612041

SITE: South Cavalcade

I. RECEIPT

- A. DATE: 12/09/86
- B. NO. OF SAMPLES: 6
- C. SAMPLE TYPE: WATER
- D. ANALYSIS REQUESTED: Full organics following EPA CLP protocol
- E. SHIPPING PROBLEMS: NONE
- F. DOCUMENTATION PROBLEMS: NONE
- G. QC DISTRIBUTION: _____

LEVEL	MATRIX	VOA	SV	PEST
LOW	WATER			
	SOIL			
MEDIUM	WATER			
	SOIL			

II. PREPARATION

- A. ANALYTICAL LEVELS: Low level

- B. EXTRACTION PROBLEMS: None

III. ANALYSIS

- A. ANALYSIS PROBLEMS: Most samples were diluted due to the high concentration of organics.

- B. SURROGATE RECOVERIES: Most surrogates were diluted out. The samples without dilution had surrogate recoveries within the range of method requirement.

- IV. OTHER PROBLEMS None

Thomas J. Oltman
Project Manager

SPECTRIX-HOUSTON

NARRATIVE - INORGANIC ANALYSES

CLIENT: Spectrix-Monroeville
LAB NO: 86-12-041

SITE: South Cavalcade

I. RECEIPT

- A. DATE: 12/09/86
B. NO. OF SAMPLES: 6
C. SAMPLE TYPE: Water
D. ANALYSIS REQUESTED: Metals: Ag, As, Be, Cd, Cr, Cu, Fe, Hg, Ni, Pb,
Sb, Se, Tl, Zn and CN

II. PREPARATION

A. REFERENCE METHOD: EPA CLP Protocol (SOW784)

B. ANALYTICAL LEVELS: Low level

C. DIGESTION PROBLEMS: None

III. ANALYSIS

A. REFERENCE METHOD: EPA CLP Protocol (SOW784)

B. ANALYSIS PROBLEMS: None

C. SPIKE RECOVERIES: None

IV. OTHER PROBLEMS None

Ken W. Euse
Department Supervisor

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INTRODUCTION TO THE SPECTRIX ORGANIC DATA PACKAGE

The purpose of this introduction is to clarify several technical report sheets contained in this report. Data sheets showing analytes and concentration are called Organic Analysis Data Sheets (OADS). OADS pages 1 and 2 contain Hazardous Substance List (HSL) compounds associated with Volatile Organics and Semi-Volatile Organics (Base/Neutral/Acids). OADS pages 4 and 5 contain Tentatively Identified Compounds (TIC) found in the Volatiles and Semi-Volatiles analyses. All concentrations listed on pages 1 and 2 are stated in UG/L or UG/KG. Header information listed on pages 1 and 2 includes client number, Spectrix internal sample number, dilution factor used, instrument used in the analysis, date received/date analyzed and other pertinent information. Soil OADS also contain listings for percent moisture and soil pH determined during the extraction process. The reporting section for pages 1 and 2 of the OADS lists individual HSL compounds and four possible listings of the concentrations found. These four possibilities are the actual concentration found, nothing detected at the detection limit as noted with a "U", a value detected but just below the detection limit noted with a "J" and a value also found in the blank noted with a "B". It is possible that a sample may contain a value just below the detection limit and also found in the blank. In this instance a "JB" would be seen after a listed detection limit. A "J" value reported indicates a trace level was found below a concentration where the quantitated value may be in question.

OADS pages 4 and 5 in general list compounds not on the HSL but also found in the GC/MS analysis. These compounds or "non-targeted compounds" are identified by comparing the peak spectra in question to a list of several thousands library spectra stored on the GC/MS magnetic disc. A measure to which the library spectra and the sample spectra are similar is the "purity". Purity is listed on the OADS as a number between 1 and 1000 where 1000 would be an extremely confident identification. OADS pages 4 and 5 list two sections: a header and data reporting section. The header contains client sample name, instrument ID, analyst and data file name. The data section shows cas #, compound name, scan #, purity of each compound and UG/L or UG/KG found. The cas #, or chemical abstract service number, is an identifying number related to the organic compound name and can help clarify differences between common names and IUPAC terminology. The scan # relates to the location of the compound in the run (i.e., retention time). The concentration listed is estimated based on the Internal Standard concentrated which is added to the sample prior to analysis. These estimated values are listed as a "J" and are not to be confused with "J" listed on OADS pages 1 and 2. Only purity values greater than 750 will show the name and cas #. Purity values less than 750 are considered unknowns and only a scan # and concentration are listed. Unknown compounds on the OADS pages 4 and 5 will also be listed on individual library search spectra sheets included in the data package. These raw data sheets show the top 4 matches to the unknown compounds so that best estimation of the compound's identity can be ascertained by you the data user. Additional manual identifications may be achieved at additional costs.

Quantitation report sheets are also found in this package. These data sheets depict raw header and analysis data important for record keeping purposes. The quantitation report also lists surrogate recoveries showing the % recovery of similar HSL compounds for Quality Control determinations. EPA % recovery acceptance levels are available upon request.

INTRODUCTION TO THE SPECTRIX INORGANIC DATA PACKAGE

The inorganic portion of the data package is designed to serve two purposes, namely, explain the reporting format and describe the methods used in analysis of samples.

Inorganic Analytical Report

Metals are reported in micrograms per liter (ug/L) for aqueous samples and milligrams per kilogram dry weight (mg/Kg dry wt.) for non-aqueous samples. Non-aqueous samples have been calculated to their dry weight basis. Default units for each parameter are shown just below the individual test code. In situations where the reported units differs from the default unit, the reported units will be shown next to the reported value.

Example: The default units for BE_S is mg/Kg dry wt., which indicates that report will be in dry weight basis. Report that shows mg/Kg below a numerical value means that sample was not calculated to dry weight basis.

Percent solid of non-aqueous samples, if performed, will be reported as PSOLID in %.

Reporting for non-metallic elements follow same pattern as metals; ug/L or mg/L for aqueous samples and mg/Kg for non-aqueous samples. Here too, mg/Kg dry weight indicates that samples have been corrected for dry weight.

Methodology for Analysis of Inorganic Parameters

The analysis of metals in aqueous samples is performed according to EPA Contract Laboratory Program (CLP) Protocol Statement of Work (SOW) No. 785, "Inorganic Analysis of Multi-Media, Multi-Concentration Samples", July, 1985. Silver, beryllium, cadmium, chromium, copper, iron, nickel, and zinc are analyzed by ICP emission spectrometry. Antimony, arsenic, lead, selenium and thallium are determined by furnace - AA, and mercury is determined by manual cold vapor. All furnace - AA and ICP analyses are background corrected. L'vov platformed graphite tubes are used for all furnace - AA analyses and matrix modified where applicable.

The analysis of non-aqueous samples is done according to EPA CLP Protocol SOW No. 785. This is a modified version of methods in EPA manual "Test methods for Evaluating Solid Waste", EPA SW-846, July, 1982.

The EP Toxicity leaching procedure is performed according to methods published by EPA in "Test Methods for Evaluating Solid Wastes", EPA SW-846, July, 1982. Analysis of the leachates for metals is performed by the method for aqueous samples cited above.

54300

CHAIN OF CUSTODY RECORD

Client/Project Name <i>Kopans Company, Inc.</i>			Project Location <i>South Cavalcade, Houston</i>											
Project No. <i>1573-73</i>		Field Logbook No. <i>KER SCOTT BARNES NO. 2</i>		Chain of Custody Tape No.										
Sampler: (Signature) <i>Brian H. Martin, Scott Barnes</i>														
Sample No./Identification	Date	Time	Lab Sample Number	Type of Sample	Acid/Base	Pesticide	PCBs	Asbestos	Organic Solvents	Flame Test	Plastics	Plants	Soil	REMARKS
CAV-DW11-001	86/12/09	1350	ground water	water	✓	✓	✓	✓	✓					6 bottles
CAV-DW10-001	86/12/09	1340	"	water	✓	✓	✓	✓	✓					6 bottles
Relinquished by: (Signature) <i>Susan R. Craig</i>			Date 86/12/09	Time 1815	Received by: (Signature)			Date	Time					
Relinquished by: (Signature)			Date	Time	Received by: (Signature)			Date	Time					
Relinquished by: (Signature)			Date	Time	Received for Laboratory: (Signature) <i>Mary F. Cruz</i>			Date 86/12/09	Time 18:15					
Sample Disposal Method:			Disposed of by: (Signature) <i>Mary F. Cruz</i>			Date	Time							
SAMPLE COLLECTOR Environmental Research and Technology, Inc. 696 Virginia Road Concord, MA 01742 617-369-8910 <i>Keystone Environmental Resources, Inc. 440 College Park Boston, MA 02114</i>			ANALYTICAL LABORATORY <i>Spectrix Corp 3911 Fondren, Suite 100 Houston, TX 77063</i>						ERT					
									No. 7451					

943300
CHAIN OF CUSTODY RECORD

Client/Project Name <i>Koppers Company, Inc.</i>		Project Location <i>South Cavalcade, Houston</i>						ANALYSES		
Project No. <i>1573-23</i>		Field Logbook No. <i>KER Scott Barnes No. 2</i>								
Sampler: (Signature) <i>Scott Barnes</i>		Chain of Custody Tape No. <i>N.A.</i>								
Sample No./ Identification	Date	Time	Lab Sample Number	Type of Sample	(2)	✓	✓	✓	✓	REMARKS
SCK R03-001	86/12/09	1615		ground water	②	✓	✓	✓	✓	6 bottles
SCK MW17-001	"	1630		"	1	✓	✓	✓	✓	5 bottles
Relinquished by: (Signature) <i>Sharon K. Cruz</i>				Date 86/12/09	Time 1815	Received by: (Signature)		Date	Time	
Relinquished by: (Signature)				Date	Time	Received by: (Signature)		Date	Time	
Relinquished by: (Signature)				Date	Time	Received for Laboratory: (Signature) <i>Mary F. Cruz</i>		Date 86/12/09	Time 18:15	
Sample Disposal Method:				Disposed of by: (Signature) <i>Mary F. Cruz</i>		Date	Time			
SAMPLE COLLECTOR <i>Keely Rose</i> Environmental Research and Technology, Inc. 696 Virginia Road Concord MA 01742 617-369-8910		ANALYTICAL LABORATORY <i>Spectra Corp.</i> 3911 Foxden, Suite 100 Houston, TX 77063		ERT						
				No. 1464						

1974-3-84

L T S E O C

CHAIN OF CUSTODY RECORD

Client/Project Name <i>Koppers Company, Inc.</i>		Project Location <i>South Cardcads, Houston</i>										
Project No. <i>1573-73</i>		Field Logbook No. <i>TER Scott Barnes Co. 2</i>										
Sampler: (Signature) <i>Scott Barnes</i>		Chain of Custody Tape No. <i>N.A.</i>										
Sample No./ Identification	Date	Time	Lab Sample Number	Type of Sample	ANALYSES							REMARKS
SCK-A02-001	86/12/09	1330		groundwater	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	6 bottles x 2 vials
CAV-AW10-001	86/12/09	1340		" "								2 vials
CAV-AW11-001	86/12/09	1350		" "								2 vials
SCK-MW17-001	86/12/09	1630		" "	<input checked="" type="checkbox"/>							2 vials + 1 bottle
SCK-P03-001	86/12/09	1615		" "								2 vials
<i>OK</i> <i>SKC</i>												
Relinquished by: (Signature) <i>Sharon K. Laij</i>				Date 86/12/09	Time 1815	Received by: (Signature)				Date	Time	
Relinquished by: (Signature)				Date	Time	Received by: (Signature)				Date	Time	
Relinquished by: (Signature)				Date	Time	Received for Laboratory: (Signature) <i>MF Cruz</i>				Date 86/12/09	Time 18:15	
Sample Disposal Method:				Disposed of by: (Signature) <i>J. Cruz</i>							Date	Time
SAMPLE COLLECTOR <i>Keptone Environmental Resources</i> Environmental Research and Technology, Inc. 696 Virginia Road Concord, MA 01742 617/369-8910		ANALYTICAL LABORATORY <i>Spectrix Corp.</i> 3911 Fondren Suite 100 Houston, TX 77063								ERT		
										Nº 1465		

8-84

A

CHECKLIST

SECTION II - Sample Data Packets

Sample No: CAV-OW10-001

Organic Analysis Data Sheets (OADS)

VOA
 SV
 Pest.

Tentatively Identified Compounds

VOA
 SV

Raw Data

VOA chromatogram
 VOA quantitation report (surrogates)
 SV chromatogram
 SV quantitation report (surrogates)
 Pest. chromatogram and integrator print out

Inorganic Analysis Data Sheets (IADS)

003348

SPECTRIX DC # ----- 8

INST ID 4000

SAMPLE NUMBER: CAV-DW10-001

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

BORATORY NAME: SPECTRIX
 SAMPLE ID NO.: 861204102
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *[Signature]*

CASE NO.: -----
 GC REPORT NO.: -----
 CONTRACT NO.: -----
 DATE SAMPLE RECEIVED: 12/9/86

VOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED/PREPARED: 12/13/86
 DATE ANALYZED: 12/13/86

DATAFILE: 4U12041V02
 DILUTION FACTOR: 1:20.....

CAS #	COMPOUND	UG/L	003549
C010	CHLOROMETHANE	200 U	
C015	BROMOMETHANE	200 U	
C020	VINYL CHLORIDE	200 U	
C025	CHLOROETHANE	200 U	
C030	METHYLENE CHLORIDE	140	
C035	ACETONE	200 U	
C040	CARBON DISULFIDE	100 U	
C045	1, 1-DICHLOROETHENE	100 U	
C050	1, 1-DICHLOROETHANE	100 U	
C055	TRANS-1, 2-DICHLOROETHENE	100 U	
C060	CHLOROFORM	100 U	
C065	1, 2-DICHLOROETHANE	100 U	
C110	2-BUTANONE	200 U	
C115	1, 1, 1-TRICHLOROETHANE	100 U	
C120	CARBON TETRACHLORIDE	100 U	
C125	VINYL ACETATE	200 U	
C130	BROMODICHLOROMETHANE	100 U	
C140	1, 2-DICHLOROPROPANE	100 U	
C145	TRANS-1, 3-DICHLOROPROPENE	100 U	
C150	TRICHLOROETHENE	100 U	
C155	DIBROMOCHLOROMETHANE	100 U	
C160	1, 1, 2-TRICHLOROETHANE	100 U	
C165	BENZENE	290	
C170	CIS-1, 3-DICHLOROPROPENE	100 U	
C175	2-CHLOROETHYL VINYLETHER	200 U	
C180	BROMOFORM	100 U	
C210	2-HEXANONE	200 U	
C205	4-METHYL-2-PENTANONE	200 U	
C220	TETRACHLOROETHENE	100 U	
C225	1, 1, 2, 2-TETRACHLOROETHANE	100 U	
C230	TOLUENE	620	
C235	CHLOROBENZENE	100 U	
C240	ETHYLBENZENE	270	
C245	STYRENE	100 U	
C250	TOTAL XYLEMES	960	

- = UNDETECTED AT THE LISTED DETECTION LIMIT
- = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT
- = COMPOUND ALSO FOUND IN BLANK

-- 8
INST ID: 4530

SPECTRIX DC # ---- 8

SAMPLE NUMBER: CAV-DW10-001

ORGANICS ANALYSIS DATA SHEET - PAGE 2

LABORATORY NAME: SPECTRIX
LAB SAMPLE ID NO.: 861204102
SAMPLE MATRIX: WATER
DATA RELEASE AUTHORIZED BY: *ymh*

CASE NO.: ---
QC REPORT NO.: ---
CONTRACT NO.: ---
DATE SAMPLE RECEIVED: *12/9/86*

SEMIVOLATILES

CONCENTRATION: LOW
DATE EXTRACTED: *12/13/86*
DATE ANALYZED: 01/09/87

DATAFILE: 4U12041C02A
DILUTION FACTOR: *400*

CAS #	COMPOUND	UG/L	0 50
C315	PHENOL	4000 U	
C325	BIS(2-CHLOROETHYL)ETHER	4000 U	
C330	2-CHLOROPHENOL	4000 U	
C335	1, 3-DICHLOROBENZENE	4000 U	
C340	1, 4-DICHLOROBENZENE	4000 U	
C345	BENZYL ALCOHOL	4000 U	
C350	1, 2-DICHLOROBENZENE	4000 U	
C355	2-METHYLPHENOL	2000 J	
C360	BIS(2-CHLOROISOPROPYL)ETHER	4000 U	
C365	4-METHYLPHENOL	1400 J	
C370	N-NITROSODIPropylAMINE	4000 U	
C375	HEXAChLORoETHANE	4000 U	
C410	NITROBENZENE	4000 U	
C415	ISOPHORONE	4000 U	
C420	2-NITROPHENOL	4000 U	
C425	2, 4-DIMETHYLPHENOL	2700 J	
C430	BENZOIC ACID	20000 U	
C435	BIS(2-CHLOROETHOXY)METHANE	4000 U	
C440	2, 4-DICHLOROPHENOL	4000 U	
C445	1, 2, 4-TRICHLOROBENZENE	4000 U	
C450	NAPHTHALENE	66000	
C455	4-CHLORoANILINE	4000 U	
C460	HEXAChLORoBUTADIENE	4000 U	
C465	P-CHLORo-M-CRESOL	4000 U	
C470	2-METHyLNAPHTHALENE	6500	
C510	HEXAChLORoCYCLOPENTADIENE	4000 U	
C515	2, 4, 6-TRICHLOROPHENOL	4000 U	
C520	2, 4, 5-TRICHLOROPHENOL	20000 U	
C525	2-CHLORoNAPHTHALENE	4000 U	
C530	2-NITROANILINE	20000 U	
C535	DIMETHYL PHTHALATE	4000 U	
C540	ACENAPHTHYLENE	290 J	
C545	3-NITROANILINE	20000 U	
C550	ACENAPHTHENE	6100	
C555	2, 4-DINITROPHENOL	20000 U	
C560	4-NITROPHENOL	20000 U	
C565	DIBENZOFURAN	4700	
C570	2, 4-DINITROTOLUENE	4000 U	
C575	2, 6-DINITROTOLUENE	4000 U	

SPECTRIX DC # ---- 8

SAMPLE NUMBER: CAV-DW10-001

EMISSIONS ORGANICS ANALYSIS DATA SHEET, CONTINUED

DATAFILE: 4U12041C02A

CAS #	COMPOUND	UG/L	00351
C580	DIETHYL PHTHALATE	4000 U	
C585	4-CHLOROPHENYL PHENYL ETHER	4000 U	
C590	FLUORENE	4800	
C595	4-NITROANILINE	20000 U	
C610	4, 6-DINITRO-2-METHYLPHENOL	20000 U	
C615	N-NITROSODIPHENYLAMINE	4000 U	
C625	4-BROMOPHENYL PHENYL ETHER	4000 U	
C630	HEXACHLOROBENZENE	4000 U	
C635	PENTACHLOROPHENOL	20000 U	
C640	PHENANTHRENE	12000	
C645	ANTHRACENE	1700 J	
C650	DI-N-BUTYL PHTHALATE	4000 U	
C655	FLUORANTHENE	4300	
C715	PYRENE	3100 J	
C720	BUTYL BENZYL PHTHALATE	4000 U	
C725	3, 3'-DICHLOROBENZIDINE	8000 U	
C730	BENZO(A)ANTHRACENE	670 J	
C735	BIS(2-ETHYLHEXYL)PHTHALATE	4000 U	
C740	CHRYSENE	620 J	
C760	DI-N-OCTYL PHTHALATE	4000 U	
C765	BENZO(B)FLUORANTHENE	4000 U	
C770	BENZO(K)FLUORANTHENE	4000 U	
C775	BENZO(A)PYRENE	4000 U	
C780	INDENO(1, 2, 3-CD)PYRENE	4000 U	
C785	DIBENZO(A, H)ANTHRACENE	4000 U	
C790	BENZO(GHI)PERYLENE	4000 U	

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

Lab Name: Spectrix
Case No: KEYSTONE
Lab No.: 86-12-041-02A

DC No: KEYSTONE
SMO No: CAV-DW10-001

ORGANICS ANALYSIS DATA SHEET
(page 3)

Pesticide/PCBs

Conc: Low
Date Extracted: 12/11/86
Date Analyzed: 12/23/86
Dilution Factor: 1

GPC Cleanup Yes ✓ No
Sep. Funnel Ext. Yes
Cont. Liquid Ext. Yes

Percent Moisture(decanted):

CAS #	Compound	ug/L
319-84-6	ALPHA-BHC	0.05 U
319-85-7	BETA-BHC	0.05 U
319-86-8	DELTA-BHC	0.05 U
58-89-9	GAMMA-BHC (LINDANE)	0.05 U
76-44-8	HEPTACHLOR	0.05 U
309-00-2	ALDRIN	0.05 U
1024-57-3	HEPTACHLOR EPOXIDE	0.05 U
959-98-8	ENDOSULFAN I	0.05 U
60-57-1	DIELDRIN	0.10 U
72-55-9	4,4'-DDE	0.10 U
72-20-8	ENDRIN	0.10 U
33213-65-9	ENDOSULFAN II	0.10 U
72-54-8	4,4'-DDD	0.10 U
7421-93-4	ENDRIN ALDEHYDE	0.10 U
1031-07-8	ENDOSULFAN SULFATE	0.10 U
50-29-3	4,4'-DDT	0.10 U
72-43-5	METHOXYSYCHLOR	0.50 U
53494-70-5	ENDRIN KETONE	0.10 U
57-74-9	CHLORDANE	0.50 U
8001-35-2	TOXAPHENE	1.0 U
12674-11-2	AROCLOTR-1016	0.50 U
11104-28-2	AROCLOTR-1221	0.50 U
11141-16-5	AROCLOTR-1232	0.50 U
53469-21-9	AROCLOTR-1242	0.50 U
12672-29-6	AROCLOTR-1248	0.50 U
11097-69-1	AROCLOTR-1254	1.0 U
11096-82-5	AROCLOTR-1260	1.0 U

003352

V_s = 500 = Volume of water extracted (ml)

V_t = 5000 = Volume of total extract (ul)

V_i = 1 = Volume of extract injected(ul)

1st Column Inst. : 6000
2nd Column Inst. : 6000 N/A

U=UNDETECTED AT THE LISTED DETECTION LIMIT.

J=COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT.

B=COMPOUND ALSO FOUND IN BLANK.

INST ID: 4000

SPECTRIX DC # ----- 8

SAMPLE NUMBER: CAV-0W10-001

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: -----

REPORT NO.: -----

ANALYST: CKT

DATAFILE: 4U12041V02

B. TENTATIVELY IDENTIFIED COMPOUNDS

AS #	VOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
24-38-9	CARBON DIOXIDE (ACN) *	34	989	2300 ^{UG/L} _m
24-38-9	CARBON DIOXIDE (ACN) *	43	979	460 ^m _{ug}
				00

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

* ALSO FOUND IN BLANK

INST ID: 4530

SPECTRIX DC # ---- 8

SAMPLE: CAV-DW10-001

ORGANICS ANALYSIS DATA SHEET - PAGE 5

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: ---

QC REPORT NO.: ---

ANALYST: GU

DATAFILE: 4U12041C02A

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	SEMIVOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
108-98-3	BENZENE, 1,3-DIMETHYL-	340	947	1600 JU
171-89-6	BENZOFURAN	519	900	2700 MJ
611-15-4	BENZENE, 1-ETHENYL-2-METHYL-	570	956	3700 OJ
95-13-6	1H-INDENE	580	944	5300 J
1498-81-7	CYCLOHEXANEMETHANOL, . ALPHA., . ALPHA., 4-TR	701	814	2400 J
10-75-0	PHENOL, 2,3-DIMETHYL-	729	906	2900 J
95-15-8	BENZO[B]THIOPHENE	752	894	4800 J
3855-26-3	PHENOL, 2-ETHYL-4-METHYL-	822	835	1600 J
171-83-2	1H-INDENE, 1-ETHYLDENE-	874	908	3800 J
575-37-1	NAPHTHALENE, 1,7-DIMETHYL-	976	944	2300 J
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	*1046		2000 J
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	1212		3800 J
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	1219		2400 J
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	1253		1700 J
86-74-8	9H-CARBAZOLE	1314	889	2800 J

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

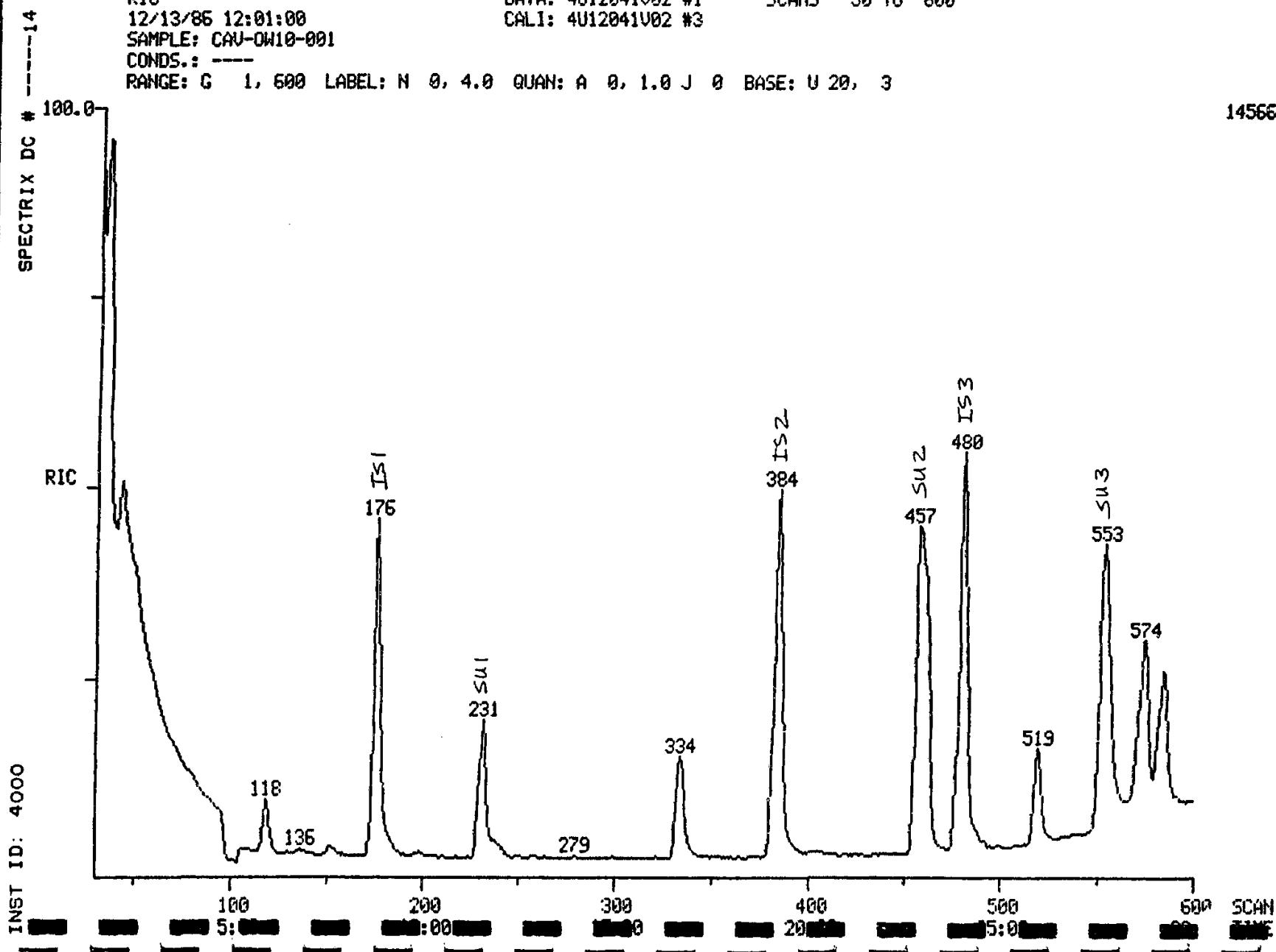
* ALSO FOUND IN BLANK

003355

RIC
12/13/86 12:01:00
SAMPLE: CAV-DW10-001
COND.: ----
RANGE: G 1, 600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12041U02 #1 SCANS 30 TO 600
CALI: 4U12041U02 #3

145664.



003355

SPECTRIX DC # -----14

VOA QUANTITATION REPORT

SAMPLE ID.	CAV-QW10-001		
ENAME	4U12041V02	INST	4000
CLIENT	KEYSTONE	ANALYST	CKT
DATE INJECTED	12/13/86 12:01:00	VERIFIED BY	<i>YDH</i>
ALIB STD DATE	121486	CORR. FACTOR	20.00

CAN#	VOA COMPOUNDS	M/E	AMOUNT	RRT	AREA
176 C101	BROMOCHLOROMETHANE	**IS1**	128	1000 UG/L	1.000 27478
384 C110	1, 4-DIFLUOROBENZENE	**IS2**	114	1000 UG/L	1.000 118229
480 C120	D5-CHLOROBENZENE	**IS3**	117	1000 UG/L	1.000 94645
231 CS15	D4-1, 2-DICHLOROETHANE	**SU1**	65	93 %	1.313 26318
457 CS05	DB-TOLUENE	**SU2**	98	97 %	0.952 76362
553 CS10	P-BROMOFLUOROBENZENE	**SU3**	95	99 %	1.152 48485
118 C030	METHYLENE CHLORIDE *		84	196 UG/L	0.670 6174
334 C165	BENZENE		78	292 UG/L	0.870 31309
461 C230	TOLUENE		92	622 UG/L	0.960 38461
519 C240	ETHYLBENZENE		106	270 UG/L	1.081 86620
574 C250	TOTAL XYLEMES		106	958 UG/L	1.196 380620

* BLANK CORRECTED ON DATA SHEET

L S E O O

RIC

01/09/87 17:49:00
SAMPLE: CAV-OW10-001
COND.: ---

DATA: 4U12041C02A #1
CALI: 4U12041C02A #3
SCANS 200 TO 1350

RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

269824

SPECTRIX DC # ----14

INST ID: 4530

100.0

RIC

200

3:

340

435

537

580

752

858

1030

753

1271

754

1118

1143

1219

1314

1000

16:40

1200

20:00

SCAN
TIME

003357

853300

RIC
01/09/87 17:49:00
SAMPLE: CAV-0W10-001
COND.: ---
RANGE: G 1,2400 LABEL: N 0, 4.0

DATA: 4U12041C02A #1 SCANS 1350 TO 2400
CALI: 4U12041C02A #3

QUAN: A 0, 1.0 J 0 BASE: U 20, 3

34816.

SPECTRIX DC # -----14

100.0

RIC

INST ID: 4530

1400
3:20

1600
26:40

1800
30:00

2000
33:20

2200
36:40

2400 SCAN
00 TIME

1409

1505

1556

1643

1710

1755

1876

1930

1936

SPECTRIX DC # ----14

SEMIVOLATILE WATER REPORT

SAMPLE ID.

CAV-DW10-001

FILENAME

4U12041C02A

INST ID.

4530

CLIENT

KEYSTONE

ANALYST

GU

DATE INJECTED

01/09/87 17:49:00

VERIFIED BY

CALIB STD DATE

01/08/87

CORR. FACTOR

400.00

SCAN#		SEMIVOLATILE COMPOUNDS	M/E	AMOUNT	RRT	AREA
536	CI30	1, 4-DICHLOROBENZEND-D4 ** IS1 **	152	16000 UG/L	1.000	36004
741	CI40	NAPHTHALENE-D8 ** IS2 **	136	16000 UG/L	1.000	136914
1030	CI50	ACENAPHTHENE-D10 ** IS3 **	164	16000 UG/L	1.000	70564
1271	CI60	PHENANTHRENE-D10 ** IS4 **	188	16000 UG/L	1.000	90041
710	CI70	CHRYSENE-D12 ** IS5 **	240	16000 UG/L	1.000	47080
1930	CI75	PERYLENE-D12 ** IS6 **	264	16000 UG/L	1.000	25171
599	C355	2-METHYLPHENOL	108	2167 UG/L	1.118	74564
623	C365	4-METHYLPHENOL	108	1409 UG/L	1.162	4833
708	C425	2, 4-DIMETHYLPHENOL	122	2705 UG/L	0.955	8441
746	C450	NAPHTHALENE	128	66389 UG/L	1.007	54565
858	C470	2-METHYLNAPHTHALENE	142	6493 UG/L	1.158	49546
1002	C540	ACENAPHTHYLENE	152	289 UG/L	0.973	2037
1035	C550	ACENAPHTHENE	153	6127 UG/L	1.005	24837
1061	C565	DIBENZOFURAN	168	4651 UG/L	1.030	25564
124	C580	DIETHYL PHTHALATE *	149	252 UG/L	1.091	1072
118	C590	FLUORENE	166	4835 UG/L	1.085	22257
1275	C640	PHENANTHRENE	178	12066 UG/L	1.003	65361
1281	C645	ANTHRACENE	178	1702 UG/L	1.008	9361
1470	C655	FLUORANTHENE	202	4309 UG/L	1.157	21837
1505	C715	PYRENE	202	3064 UG/L	0.880	13841
1708	C730	BENZO(A)ANTHRACENE	228	665 UG/L	0.999	2036
1713	C740	CHRYSENE	228	616 UG/L	1.002	1751

* BLANK CORRECTED ON DATA SHEET

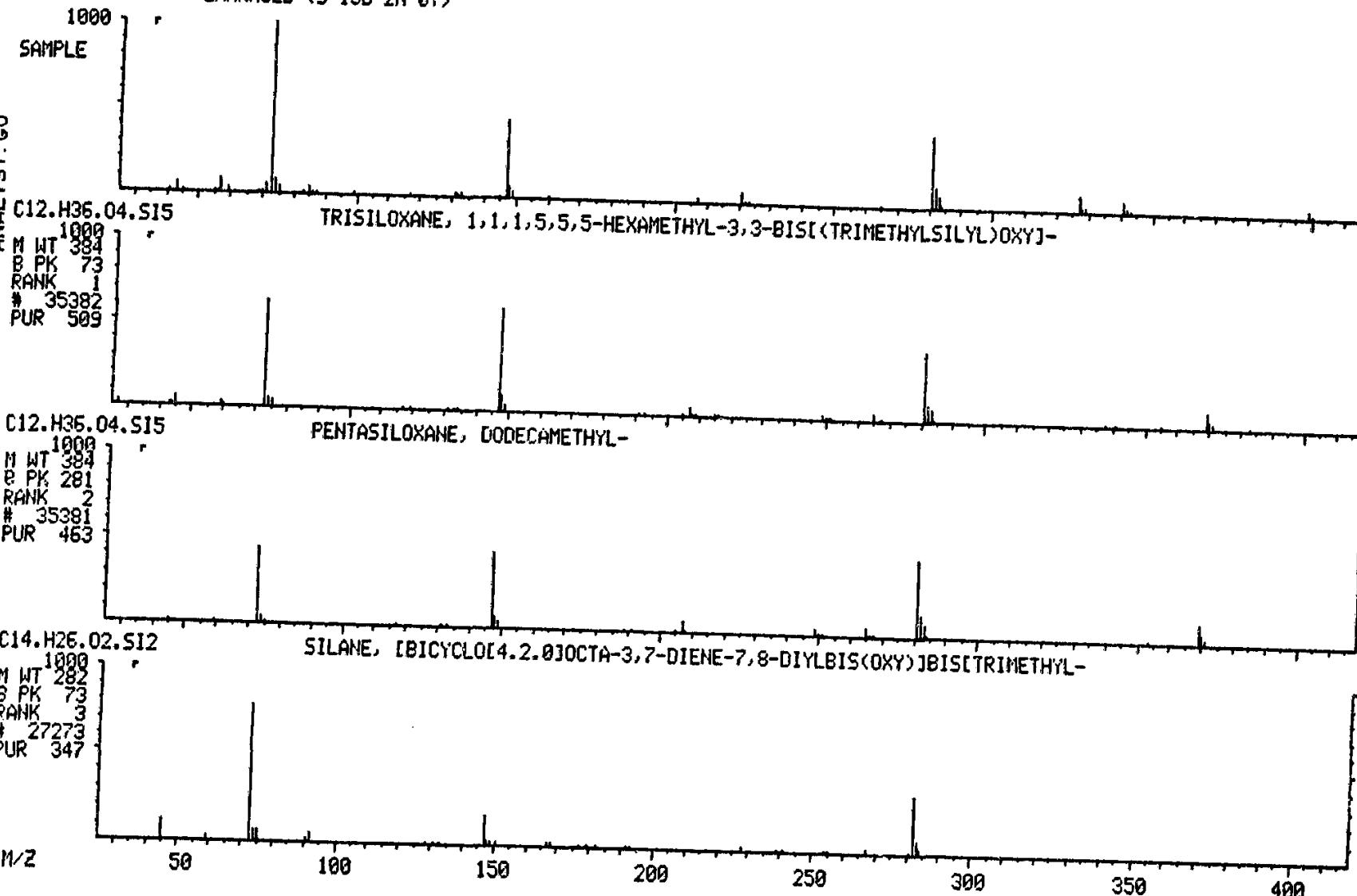
** SURROGATES WERE DILUTED OUT

09 E E 00

LIBRARY SEARCH
01/09/87 17:49:00 + 17:26
SAMPLE: CAV-0W10-001
COND.: ---
ENHANCED (S 15B 2N 0T)

DATA: 4U12041C02A #1046 BASE M/Z: 73
CALI: 4U12041C02A # 3 RIC: 8575.

SPECTRIX DC # -----19
ANALYST: GU



193300

LIBRARY SEARCH
01/09/87 17:49:00 + 20:12
SAMPLE: CAV-0W10-001
COND.: ---
ENHANCED (S 15B 2N 0T)

DATA: 4U12041C02A #1212 BASE M/Z: 170
CALI: 4U12041C02A # 3 RIC: 3159.

SPECTRUM DC # -----19

1263
SAMPLE

C14.H12
M WT 1263
B PK 165
RANK 1
13177
PUR 427

9H-FLUORENE, 4-METHYL-

C14.H12
M WT 1263
B PK 165
RANK 2
13179
PUR 424

9H-FLUORENE, 9-METHYL-

C12.H10.0
M WT 1263
B PK 170
RANK 3
11441
PUR 412

[1,1'-BIPHENYL]-3-OL

INST ID: 4530

M/Z 40 60 80 100 120 140 160 180 200 220

003360

LIBRARY SEARCH
01/09/87 17:49:00 + 20:19
SAMPLE: CAV-0W10-001
COND.: ---
ENHANCED (S 15B 2N 0T)

DATA: 4U12041C02A #1219 BASE M/Z: 145
CALI: 4U12041C02A # 3 RIC: 4847.

SPECTRUM DC # -----19

1348
SAMPLE

C9.H7.O.N
M WT 1348
B PK 145
RANK 1
6784
PUR 479

5-ISOQUINOLINOL

C9.H7.O.N
M WT 1348
B PK 145
RANK 2
6772
PUR 475

8-QUINOLINOL

C9.H7.O.N
M WT 1348
B PK 145
RANK 3
6783
PUR 476

QUINOLINE, 1-OXIDE

M/Z

40 60 80 100 120 140 160 180

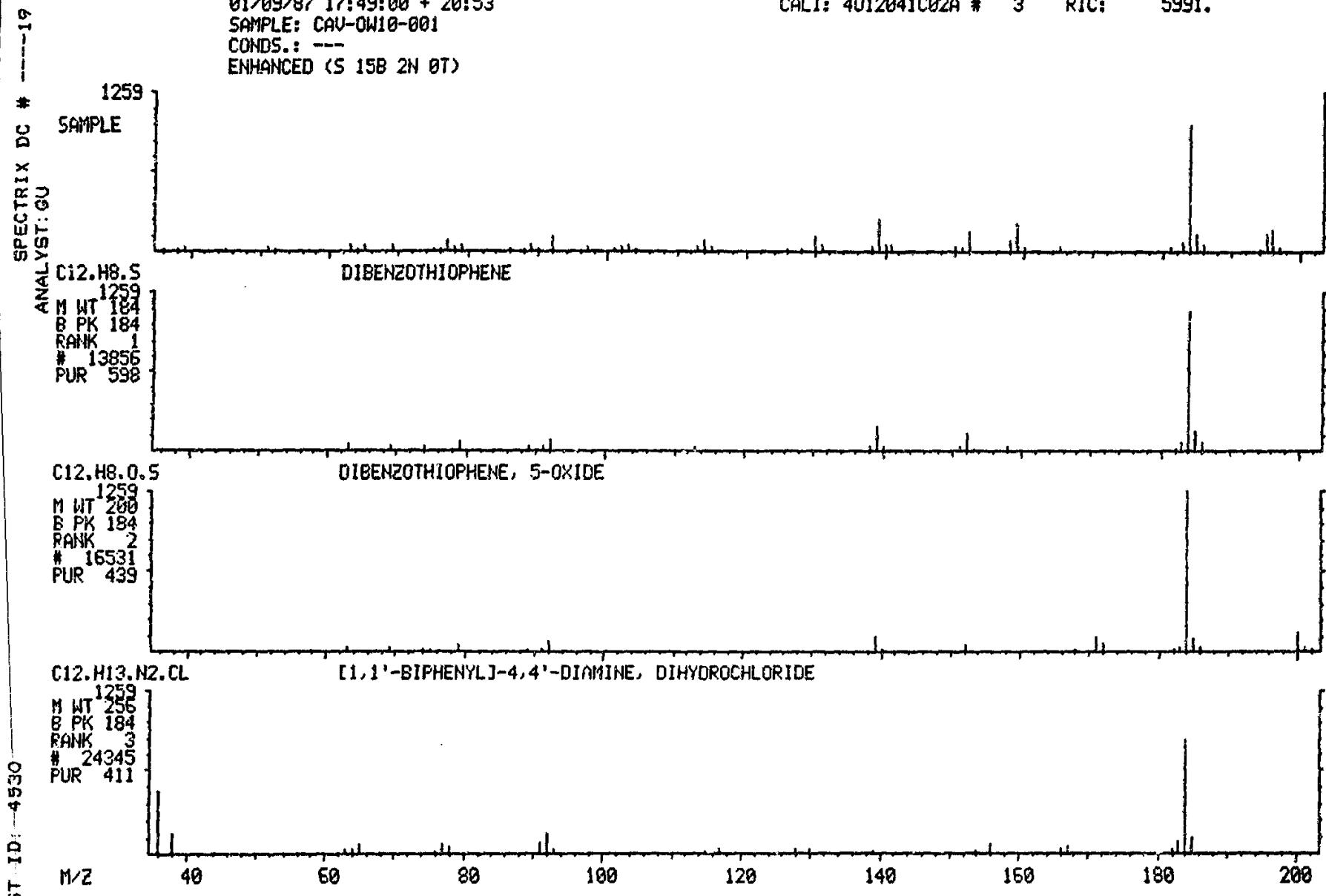
INST ID: 4530

003362

595300

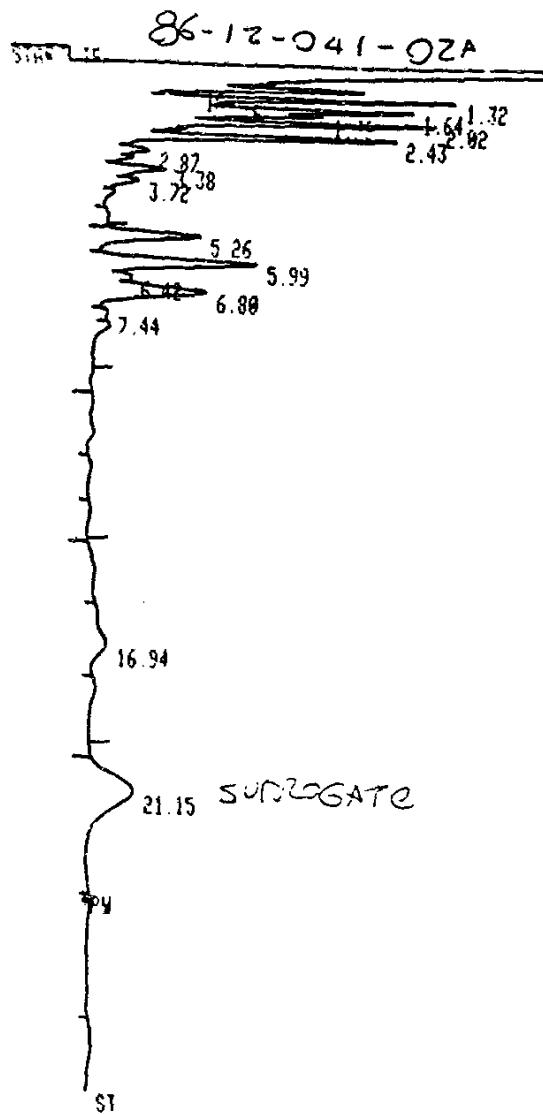
LIBRARY SEARCH
01/09/87 17:49:00 + 20:53
SAMPLE: CAV-0W10-001
COND.: ---
ENHANCED (S 15B 2N 0T)

DATA: 4U12041C02A #1253 BASE M/Z: 184
CALI: 4U12041C02A # 3 RIC: 5991.



INST ID: 4530

003363



DC # N/A

SPECTRIX CORPORATION

CASE NO. Keystone

SAMPLE NO. CAV-SW10-001

PESTICIDE FRACTION

1st Column - 1D

1.8mm x 4mm ID x 6mm 00

1.5% SP 2250

1.95% SP2401

100/120 Supelcoport

Column Temp. 203 °C

Volume Inj. 1.0 ul

Instrument ID VARIAN 6c

RUN # 135
WORKFILE ID: C
WORKFILE NAME:

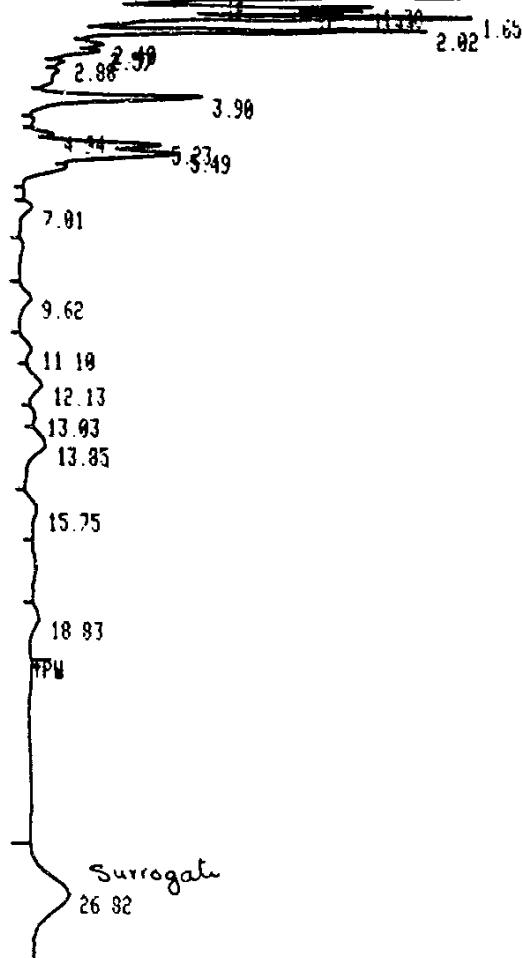
HEIGHT:

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.32	38415	D BY	0.105	16.001
1.64	33876	VV	0.135	14.111
1.72	22534	D VV	0.130	9.386
2.02	106	D VV	0.123	15.273
2.43		VV	0.129	14.086
2.87	204	VV	0.153	1.352
3.38	6423	PY	0.192	2.675
3.72	3253	VP	0.175	1.355
5.26	12403	SP	0.240	5.166
5.99	20386	PV	0.260	8.492
6.42	4297	VV	0.225	1.798
6.88	14175	VV	0.346	5.904
7.44	1559	VV	0.290	0.649
16.94	2187	VV	1.300	0.911
21.15	5636	BP	1.106	2.348

TOTAL HGT= 240020
MULT. FACTOR= 1.0000E+00

STAR 86-12-041-02A

NA.



DC#
SPECTRIX HOUSTON

CASE NO. Keystone

SAMPLE NO. CAV-0W10-0
PESTICIDE FRACTION
2ND COLUMN CONFIRMATI

2ND COLUMN II C
1.8m x2mm ID x6mm OD
3% SP-2100
100/120 SUPELCOPORT
COLUMN TEMP. 187 °C
VOLUME INJ. 1 μl
INSTRUMENT ID. 6000

M
O
O

RUN # 12 JAN/06/82 17:49:24
WORKFILE ID: C
WORKFILE NAME:

HEIGHT%

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.38	61678	BV	0.097	11.865
1.49	60773	D VV	0.099	11.691
1.65	134658	D VP	0.100	25.902
2.02	85865	D PV	0.104	16.518
2.40	7287	VV	0.161	1.402
2.57	8483	D VP	0.146	1.632
2.98	2323	D PV	0.102	0.447
3.90	42327	PP	0.197	8.142
4.94	6088	VV	0.198	1.171
5.23	33659	VV	0.209	6.475
5.49	38786	VV	0.272	7.446
7.01	2678	VP	0.320	0.515
9.62	3169	PP	0.551	0.610
11.10	2385	PV	0.532	0.555
12.13	5241	VV	0.765	1.088
13.03	3457	VV	0.527	0.665
13.85	5627	VV	0.851	1.083
15.75	2957	VV	0.723	0.569
18.83	2429	VB	0.687	0.467
26.82	9554	BB	1.243	1.838

Page 1
Received: 12/09/86

SPECTROCHEMICAL REPORT
Results by Sample

Work Order # 86-12-041

SAMPLE ID CAV-DW10-001 (LLW)		SAMPLE # 02 FRACTIONS: A			
Date & Time Collected 12/09/86			Category _____		
AG	<10	AS	34.8	BE	<3
	ug/L		ug/L		ug/L
CU	33	FE	35500	HG	<0.2
	ug/L		ug/L		ug/L
SB	67	SE	<5	TL	<10
	ug/L		ug/L		ug/L
				ZN	116
					ug/L

CHECKLIST

SECTION II - Sample Data Packets

Sample No: CAV-OWII-001

✓ Organic Analysis Data Sheets (OADS)

✓ VOA
✓ SV
✓ Pest.

✓ Tentatively Identified Compounds

✓ VOA
✓ SV

✓ Raw Data

✓ VOA chromatogram
✓ VOA quantitation report (surrogates)
✓ SV chromatogram
✓ SV quantitation report (surrogates)
✓ Pest. chromatogram and integrator print out

✓ Inorganic Analysis Data Sheets (IADS)

003367

SPECTRIX DC # ----- 8

INST ID 4000

SAMPLE NUMBER: CAV-DW11-001

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

BORATORY NAME: SPECTRIX
 LMB SAMPLE ID NO.: 861204103
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *[Signature]*

CASE NO.: -----
 GC REPORT NO.: -----
 CONTRACT NO.: -----
 DATE SAMPLE RECEIVED: 12/13/86

VOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED/PREPARED: 12/13/86
 DATE ANALYZED: 12/13/86

DATAFILE: 4U12041V03

DILUTION FACTOR: 1/100....

CAS #	COMPOUND	UG/L
C010	CHLOROMETHANE	1000 U
C015	BROMOMETHANE	1000 U
C020	VINYL CHLORIDE	1000 U
C025	CHLOROETHANE	1000 U
C030	METHYLENE CHLORIDE	1000 U
C035	ACETONE	500 U
C040	CARBON DISULFIDE	1000 J
C045	1, 1-DICHLOROETHENE	500 U
C050	1, 1-DICHLOROETHANE	500 U
C055	TRANS-1, 2-DICHLOROETHENE	500 U
C060	CHLOROFORM	500 U
C065	1, 2-DICHLOROETHANE	500 U
C110	2-BUTANONE	500 U
C115	1, 1, 1-TRICHLOROETHANE	1000 U
C120	CARBON TETRACHLORIDE	500 U
C125	VINYL ACETATE	500 U
C130	BROMODICHLOROMETHANE	1000 U
C140	1, 2-DICHLOROPROPANE	500 U
C145	TRANS-1, 3-DICHLOROPROPENE	500 U
C150	TRICHLOROETHENE	500 U
C155	DIBROMOCHLOROMETHANE	500 U
C160	1, 1, 2-TRICHLOROETHANE	500 U
C165	BENZENE	500 U
C170	CIS-1, 3-DICHLOROPROPENE	500 J
C175	2-CHLOROETHYL VINYL ETHER	500 U
C180	BROMOFORM	1000 U
C210	2-HEXANONE	500 U
C205	4-METHYL-2-PENTANONE	1000 U
C220	TETRACHLOROETHENE	1000 U
C225	1, 1, 2, 2-TETRACHLOROETHANE	500 U
C230	TOLUENE	500 U
C235	CHLOROBENZENE	500 U
C240	ETHYL BENZENE	500 U
C245	STYRENE	500 U
C250	TOTAL XYLEMES	500 J

= UNDETECTED AT THE LISTED DETECTION LIMIT

= COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

= COMPOUND ALSO FOUND IN BLANK

INST ID: 4530

SPECTRIX DC # ---- S
SAMPLE NUMBER: CAV-DW11-001

ORGANICS ANALYSIS DATA SHEET - PAGE 2

LABORATORY NAME: SPECTRIX
 LAB SAMPLE ID NO.: 8612041C3
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *[Signature]*

CASE NO.: ---
 QC REPORT NO.: ---
 CONTRACT NO.: ---
 DATE SAMPLE RECEIVED: 12/9/86

SEMICVOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED: 12/13/86
 DATE ANALYZED: 01/09/87

DATAFILE: 4U12041C03
 DILUTION FACTOR: 50000

CAS #	COMPOUND	UG/L
C315	PHENOL	500000 U
C325	BIS(2-CHLOROETHYL)ETHER	500000 U
C330	2-CHLOROPHENOL	500000 U
C335	1, 3-DICHLOROBENZENE	500000 U
C340	1, 4-DICHLOROBENZENE	500000 U
C345	BENZYL ALCOHOL	500000 U
C350	1, 2-DICHLOROBENZENE	500000 U
C355	2-METHYLPHENOL	500000 U
C360	BIS(2-CHLOROISOPROPYL)ETHER	500000 U
C365	4-METHYLPHENOL	500000 U
C370	N-NITROSODIPropylAMINE	500000 U
C375	HEXAChLOROETHANE	500000 U
C410	NITROBENZENE	500000 U
C415	ISOPHORONE	500000 U
C420	2-NITROPHENOL	500000 U
C425	2, 4-DIMETHYLPHENOL	500000 U
C430	BENZOIC ACID	2500000 U
C435	BIS(2-CHLOROETHOXY)METHANE	500000 U
C440	2, 4-DICHLOROPHENOL	500000 U
C445	1, 2, 4-TRICHLOROBENZENE	500000 U
C450	NAPHTHALENE	7100000
C455	4-CHLORoANILINE	500000 U
C460	HEXAChLOROBUTADIENE	500000 U
C465	P-CHLORo-M-CRESOL	500000 U
C470	2-METHYLNAPHTHALENE	1300000
C510	HEXAChLOROCYCLOPENTADIENE	500000 U
C515	2, 4, 6-TRICHLOROPHENOL	500000 U
C520	2, 4, 5-TRICHLOROPHENOL	2500000 U
C525	2-CHLORoNAPHTHALENE	500000 U
C530	2-NITROANILINE	2500000 U
C535	DIMETHYL PHTHALATE	500000 U
C540	ACENAPHTHYLENE	50000 U
C545	3-NITROANILINE	2500000 U
C550	ACENAPHTHENE	2600000
C555	2, 4-DINITROPHENOL	2500000 U
C560	4-NITROPHENOL	2500000 U
C565	DIBENZOFURAN	1300000
C570	2, 4-DINITROTOLUENE	500000 U
C575	2, 6-DINITROTOLUENE	500000 U

SPECTRIX DC # ---- 8

SAMPLE NUMBER: CAV-DW11-001

IN VOLATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

DATAFILE: 4U12041C03

CAS #	COMPOUND	UG/L
C580	DIETHYL PHTHALATE	500000 U
C585	4-CHLOROPHENYL PHENYL ETHER	500000 U
C590	FLUORENE	1800000
C595	4-NITROANILINE	2500000 U
C610	4, 6-DINITRO-2-METHYLPHENOL	2500000 U
C615	N-NITROSODIPHENYLAMINE	500000 U
C625	4-BROMOPHENYL PHENYL ETHER	500000 U
C630	HEXACHLOROBENZENE	500000 U
C635	PENTACHLOROPHENOL	2500000 U
C640	PHENANTHRENE	4900000
C645	ANTHRACENE	550000
C650	DI-N-BUTYL PHTHALATE	500000 U
C655	FLUORANTHENE	2600000
C715	PYRENE	1900000
C720	BUTYL BENZYL PHTHALATE	500000 U
C725	3, 3'-DICHLOROBENZIDINE	1000000 U
C730	BENZO(A)ANTHRACENE	500000
C735	BIS(2-ETHYLHEXYL)PHTHALATE	500000 U
C740	CHRYSENE	420000 J
C760	DI-N-OCTYL PHTHALATE	500000 U
C765	BENZO(B)FLUORANTHENE	500000 U
C770	BENZO(K)FLUORANTHENE	500000 U
C775	BENZO(A)PYRENE	170000 J
C780	INDENO(1, 2, 3-CD)PYRENE	500000 U
C785	DIBENZO(A, H)ANTHRACENE	500000 U
C790	BENZO(GHI)PERYLENE	500000 U

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

003370

Lab Name: Spectrix
Case No: KEYSTONE
Lab No.: 86-12-041-03A

DC No: KEYSTONE
SMO No: CAV-0W11-001

ORGANICS ANALYSIS DATA SHEET
(page 3)

Pesticide/PCBs

Conc: Low
Date Extracted: 12/11/86
Date Analyzed: 12/23/86
Dilution Factor: 5
Percent Moisture(decanted):

GPC Cleanup Yes No
Sep. Funnel Ext. Yes
Cont. Liquid Ext. Yes

CAS #	Compound	ug/L
319-34-6	ALPHA-BHC	0.25 U
319-85-7	BETA-BHC	0.25 U
319-86-8	DELTA-BHC	0.25 U
58-89-9	GAMMA-BHC (LINDANE)	0.25 U
76-44-8	HEPTACHLOR	0.25 U
309-00-2	ALDRIN	0.25 U
1024-57-3	HEPTACHLOR EPOXIDE	0.25 U
959-98-8	ENDOSULFAN I	0.25 U
60-57-1	DIELDRIN	0.50 U
72-55-9	4,4'-DDE	0.50 U
72-20-8	ENDRIN	0.50 U
33213-65-9	ENDOSULFAN II	0.50 U
72-54-8	4,4'-DDD	0.50 U
7421-93-4	ENDRIN ALDEHYDE	0.50 U
1031-07-8	ENDOSULFAN SULFATE	0.50 U
50-29-3	4,4'-DDT	0.50 U
72-43-5	METHOXYSCHLOR	2.50 U
53494-70-5	ENDRIN KETONE	0.50 U
57-74-9	CHLORDANE	2.50 U
8001-35-2	TOXAPHENE	5.0 U
12674-11-2	AROCOLOR-1016	2.50 U
11104-28-2	AROCOLOR-1221	2.50 U
11141-16-5	AROCOLOR-1232	2.50 U
53469-21-9	AROCOLOR-1242	2.50 U
12672-29-6	AROCOLOR-1248	2.50 U
11097-69-1	AROCOLOR-1254	5.0 U
11096-82-5	AROCOLOR-1260	5.0 U

V_s = 500 = Volume of water extracted (ml)

V_t = 5000 = Volume of total extract (ul)

V_i = 1 = Volume of extract injected(ul)

1st Column Inst. : 6000

2nd Column Inst. : 6000

NA.

U=UNDETECTED AT THE LISTED DETECTION LIMIT.

J=COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT.

B=COMPOUND ALSO FOUND IN BLANK.

003371

INST ID: 4000

SPECTRIX DC # ----- B

SAMPLE NUMBER: CAV-DW11-001

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: -----

REPORT NO.: -----

ANALYST: TTO

DATAFILE: 4U12041V03

B. TENTATIVELY IDENTIFIED COMPOUNDS

AS #	VOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT	UGRN
124-38-9	CARBON DIOXIDE (ACN) *	43	1000	1800	457
124-38-9	CARBON DIOXIDE (ACN) *	48	979	2600	0650

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

* ALSO FOUND IN BLANK

INST ID: 4530

SPECTRIX DC # ---- 8

SAMPLE: CAV-DW11-001

ORGANICS ANALYSIS DATA SHEET - PAGE 5

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: ---

GC REPORT NO.: ---

ANALYST: GU

DATAFILE: 4U12041C03

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	SEMIVOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
611-15-4	BENZENE, 1-ETHENYL-2-METHYL-	559	956	550000 J
95-15-8	BENZO[B]THIOPHENE	750	915	230000 J
2471-83-2	1H-INDENE, 1-ETHYLIDENE-	875	915	680000 J
2-52-4	1, 1'-BIPHENYL	940	925	430000 J
939-27-5	NAPHTHALENE, 2-ETHYL-	954	892	240000 J
31-40-8	NAPHTHALENE, 2, 3-DIMETHYL-	964	927	310000 J
581-40-8	NAPHTHALENE, 2, 3-DIMETHYL-	978	938	440000 J
644-08-6	1, 1'-BIPHENYL, 4-METHYL-	1136	900	310000 J
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	1151		230000 J
7320-53-8	DIBENZOFURAN, 4-METHYL-	1162	869	290000 J
732-65-0	DIBENZOTHIOPHENE	1255	801	420000 J
74410-42-7	D-GALACTITOL, 2-(ACETYLMETHYLAMINO)-2-DEO	1316	911	230000 J
1961-96-2	1H-INDENE, 1-PHENYL-	1359	934	220000 J
1961-96-2	1H-INDENE, 1-PHENYL-	1363	924	260000 J
203-64-5	4H-CYCLOPENTA[DEF]PHENANTHRENE	1376	816	620000 J
38-84-6	11H-BENZO[A]FLUORENE	1569	877	350000 J

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

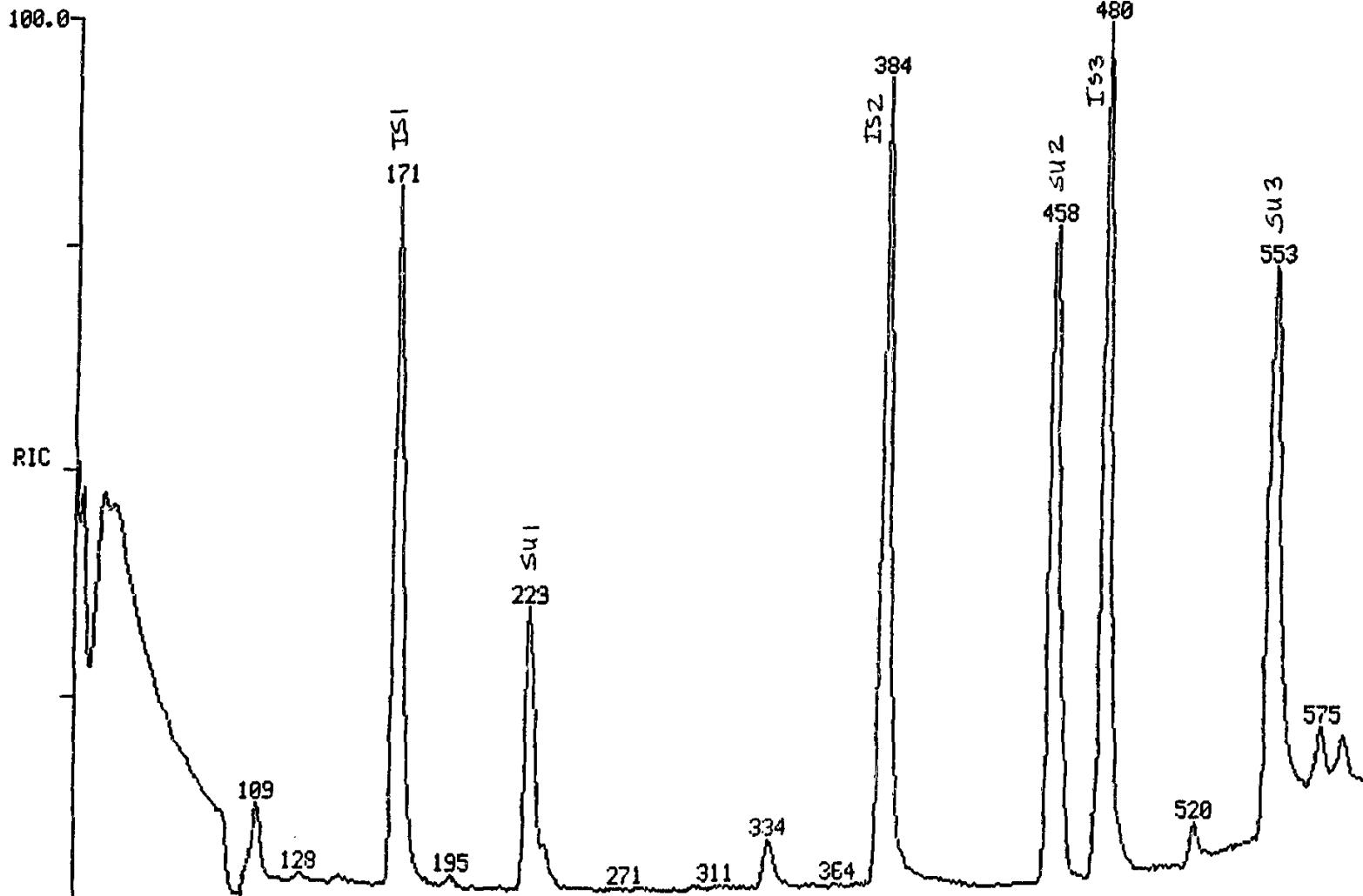
003374

RIC
12/13/86 16:54:00
SAMPLE: CAV-0W11-001
COND.: ----
RANGE: G 1, 600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12041U03 #1 SCANS 30 TO 600
CALI: 4U12041U03 #3

SPECTRIX DC # -----14

INST ID: 4000



84096.

003374

SPECTRIX DC # -----14

VOA QUANTITATION REPORT

SAMPLE ID. CAV-OW11-001
 FENAME 4U12041V03 INST 4000
 IENT KEYSTONE ANALYST TTO
 DATE INJECTED 12/13/86 16:54:00 VERIFIED BY *[Signature]*
 ALIB STD DATE 121486 CORR. FACTOR 100.00

CAN#	VOA COMPOUNDS	M/E	AMOUNT	RRT	AREA	
171 C101	BROMOCHLOROMETHANE	**IS1**	128	5000 UG/L	1.000	29148
384 C110	1, 4-DIFLUOROBENZENE	**IS2**	114	5000 UG/L	1.000	123612
480 C120	D5-CHLOROBENZENE	**IS3**	117	5000 UG/L	1.000	100136
229 CS15	D4-1, 2-DICHLOROETHANE	**SU1**	65	96 %	1.339	28626
457 CS05	D8-TOLUENE	**SU2**	98	97 %	0.952	80638
553 CS10	P-BROMOFLUOROBENZENE	**SU3**	95	96 %	1.152	50080
109 C030	METHYLENE CHLORIDE *		84	632 UG/L	0.637	4223
128 C035	ACETONE		43	545 UG/L	0.749	1327
236 C110	2-BUTANONE *		72	979 UG/L	0.615	1173
334 C165	BENZENE		78	332 UG/L	0.870	7431
461 C230	TOLUENE		92	591 UG/L	0.960	7731
520 C240	ETHYLBENZENE		106	268 UG/L	1.083	1819
584 C250	TOTAL XYLEMES		106	370 UG/L	1.217	3110

* BLANK CORRECTED ON DATA SHEET

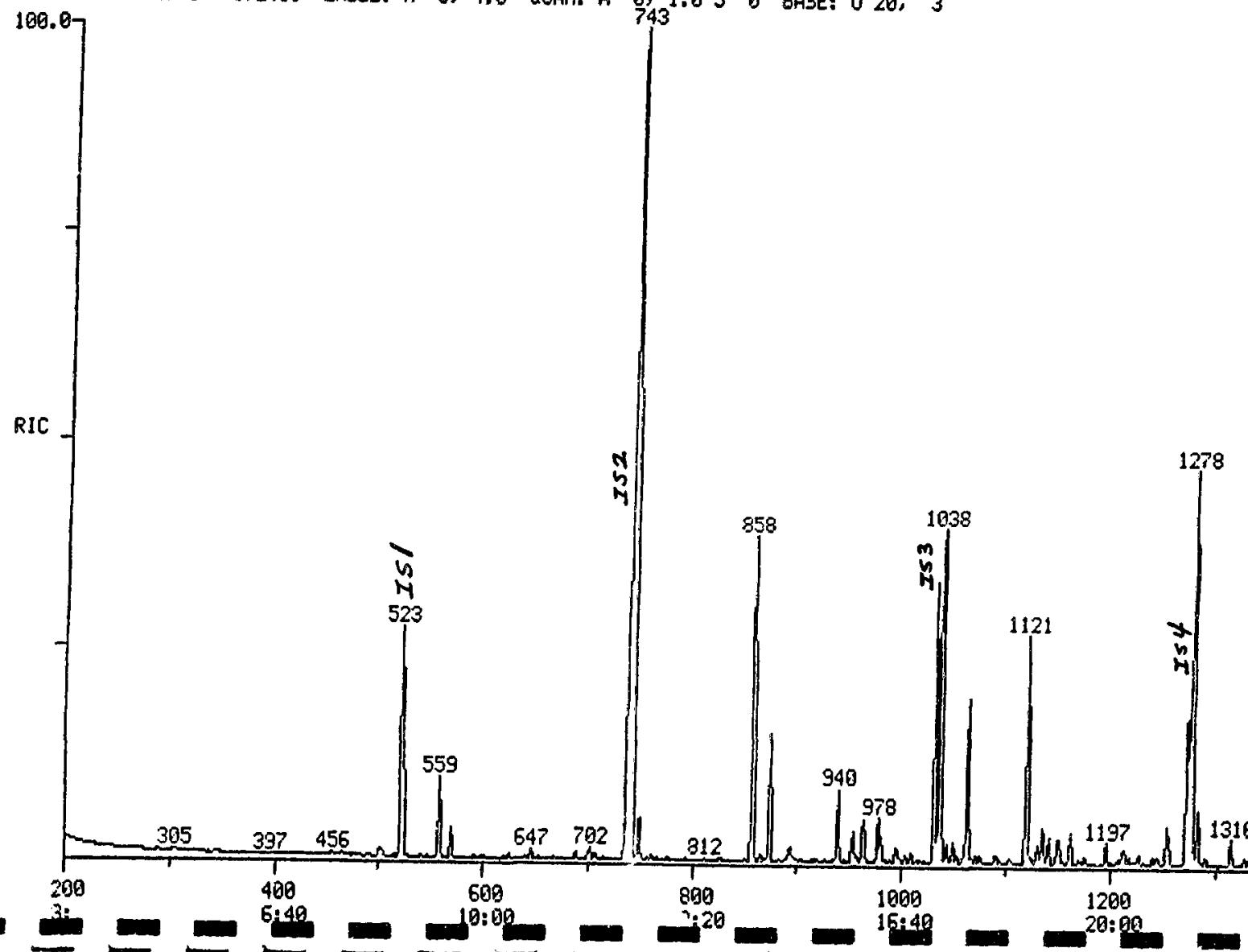
9 L E S O O

RIC
01/09/87 15:45:00
SAMPLE: CAV-OW11-001
COND.S.: ---
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12041C03 #1
CALI: 4U12041C03 #3
SCANS 200 TO 1350

SPECTRIX DC # ----14

INST ID: 4530



259072.

SCAN
TIME

LL300

RIC
01/09/87 15:45:00
SAMPLE: CAV-0W11-001
COND.: ---
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12041C03 #1 SCANS 1350 TO 2400
CALI: 4U12041C03 #3

55168.

SPECTRIX DC # ----14

100.0

RIC

1376

1412

1509

1569

1400
23:20

1600
25:40

1800
30:00

2000
33:20

2200
36:40

2400
40:00 SCAN TIME

155

1713

1932

156

1780

1878

2095

2152

2214

2283

2345

INST ID: 4530

003377

SPECTRIX DC # ----14

SEMIVOLATILE WATER REPORT

SAMPLE ID. CAV-DW11-001
 LENAME 4U12041C03
 CLIENT KEYSTONE
 DATE INJECTED 01/09/87 15:45:00
 CALIB STD DATE 01/08/87

INST ID. 4530
 ANALYST GU
 VERIFIED BY JPH
 CORR. FACTOR 50000.00

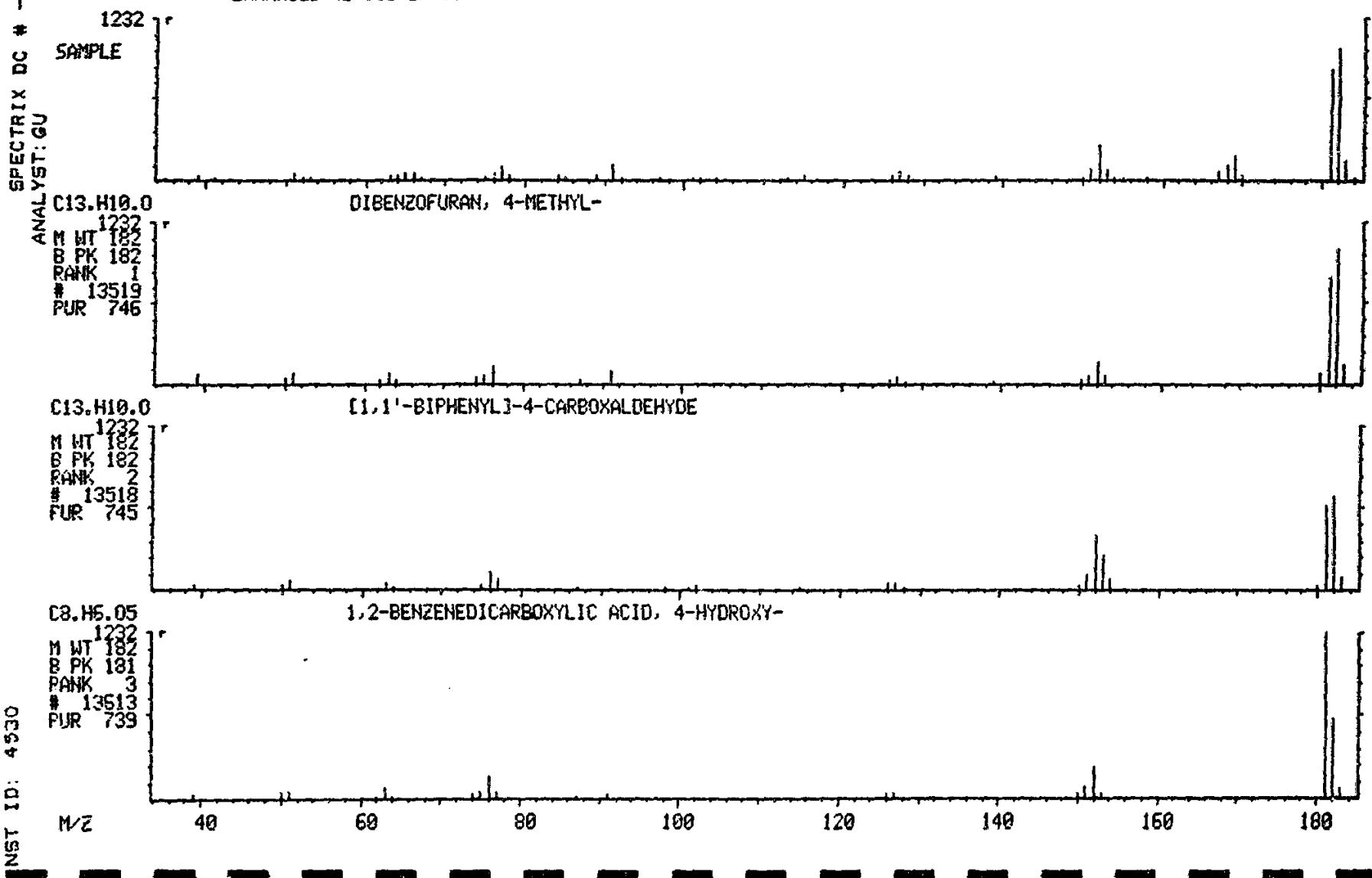
SCAN#	SEMIVOLATILE COMPOUNDS		M/E	AMOUNT	RRT	AREA
523	C130	1, 4-DICHLOROBENZEND-D4 ** IS1 **	152	2000000 UG/L	1.000	30812
738	C140	NAPHTHALENE-D8 ** IS2 **	136	2000000 UG/L	1.000	120694
1032	C150	ACENAPHTHENE-D10 ** IS3 **	164	2000000 UG/L	1.000	59246
1273	C160	PHENANTHRENE-D10 ** IS4 **	188	2000000 UG/L	1.000	60800
1713	C170	CHRYSENE-D12 ** IS5 **	240	2000000 UG/L	1.000	3467
1932	C175	PE'YLENE-D12 ** IS6 **	264	2000000 UG/L	1.000	1928
743	C450	NAPHTHALENE	128	7131280 UG/L	1.000	41333
858	C470	2-METHYLNAPHTHALENE	142	1280532 UG/L	1.163	6890
1004	C540	ACENAPHTHYLENE	152	50177 UG/L	0.973	237
1038	C550	ACENAPHTHENE	153	2608512 UG/L	1.006	7102
1064	C565	DIBENZOFURAN	168	1344266 UG/L	1.031	4964
1121	C590	FLUORENE	166	1803057 UG/L	1.086	5577
1278	C640	PHENANTHRENE	178	4900695 UG/L	1.004	15951
1284	C645	ANTHRACENE	178	548172 UG/L	1.009	1812
174	C655	FLUORANTHENE	202	2595579 UG/L	1.158	7900
509	C715	PYRENE	202	1949747 UG/L	0.881	5191
1710	C730	BENZO(A)ANTHRACENE	228	496186 UG/L	0.998	895
1716	C740	CHRYSENE	228	416407 UG/L	1.002	700
1923	C775	BENZO(A)PYRENE	252	165242 UG/L	0.995	165

* SURROGATES WERE DILUTED OUT

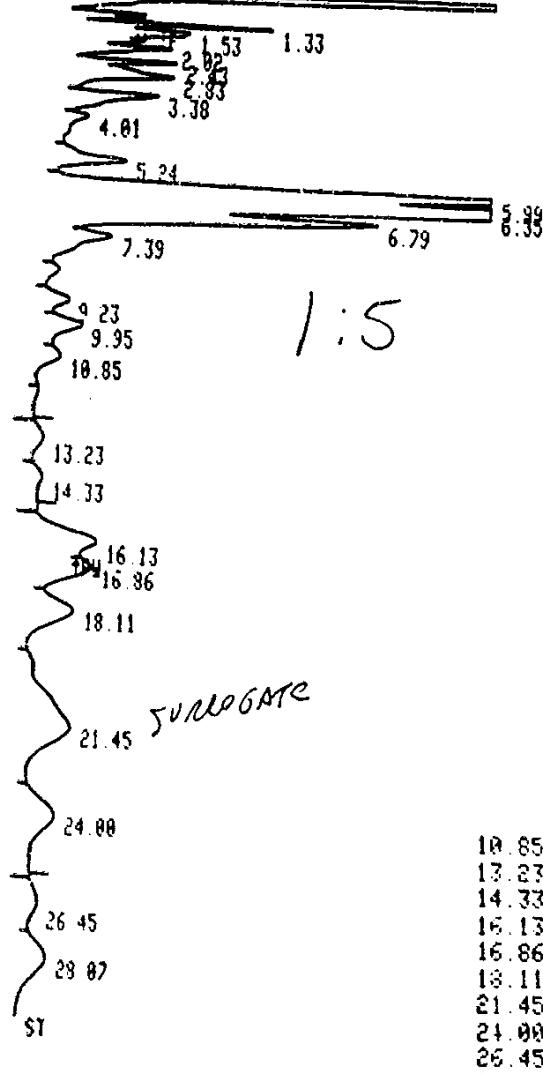
6 LΣΣ00

LIBRARY SEARCH
01/09/87 15:45:00 + 19:11
SAMPLE: CAV-0W11-001
COND.: --
ENHANCED (S 15B 2N BT)

DATA: 4U12041C03 #1151 BASE M/Z: 192
CALI: 4U12041C03 # 3 RIC: 6871.



STAR 86-12-241-03A 1:5 ROG



DC # N/D

SPECTRIX CORPORATION

CASE NO. Keytene

SAMPLE NO.CAV - 0WII-001

1:5

PESTICIDE FRACTION

1st Column - 1D

1.8m X 4mm ID X 6mm 00

0

1.5% SP 2250

00

1.95% SP2401

00

100/120 Supelcoport

00

Column Temp. 203 °C

Volume Inj. 1.0 ul

Instrument ID VARIAN 60

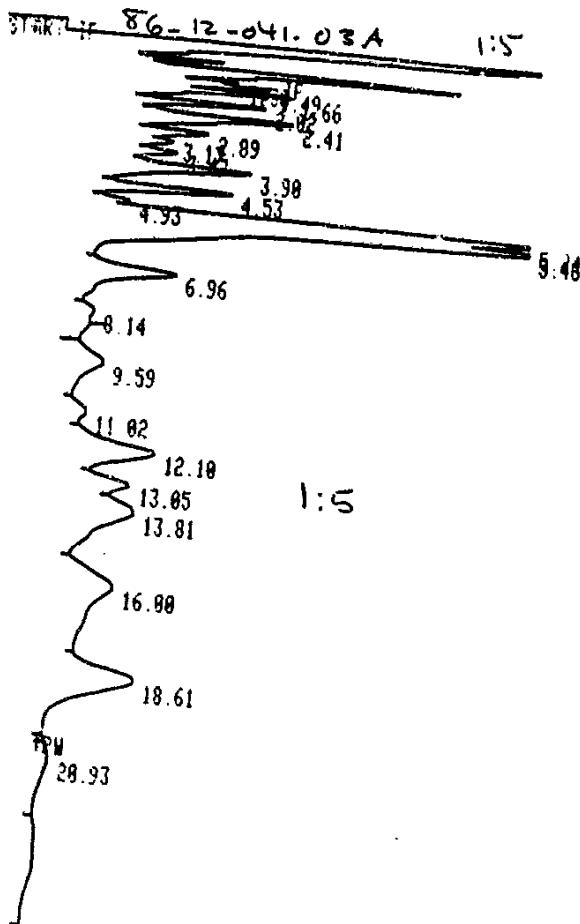
18.85	2838	VP	0.584	0.621
17.23	1318	BP	0.599	0.288
14.33	966	PB	0.492	0.211
16.13	7983	BV	0.689	1.746
16.86	7483 D	VV	0.717	1.637
18.11	5540	VP	0.805	1.212
21.45	5468	PP	1.512	1.194
24.00	3322	PE	0.998	0.727
26.45	1563	BV	0.996	0.342
28.07	3341 1	VF	1.306	0.731

RUN # 136
WORKFILE ID: C
WORKFILE NAME:

DEC/23/86 24 47:49

HEIGHT%

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.33	21917	D PV	0.093	4.794
1.53	11651	VV	0.261	2.548
2.02	10489	D VP	0.178	2.294
2.43	11972	PV	0.127	2.619
2.93	12090	VV	0.288	2.644
3.39	10738	VV	0.270	2.349
4.01	2509	VP	0.282	0.549
5.24	8589	PP	0.284	1.879
5.99	77660	PV	0.247	16.985
6.35	191209	VV	0.258	41.820
6.79	41916	VV	0.304	9.168
7.39	7971	VV	0.481	1.743
9.23	3355	VV	0.417	0.234
9.95	5337	VV	0.515	1.167



NN.

DCP
SPECTRIX HOUSTON
CASE NO. Keystone
SAMPLE NO. CAV-0WII-0c
PESTICIDE FRACTION
2ND COLUMN CONFIRMATI.
2ND COLUMN II C
1.8m x2mm ID x6mm OD
3% SP-2100
100/120 SUPELCOPORT
COLUMN TEMP. 187 °C
VOLUME INJ. 1 AL
INSTRUMENT ID. 6008

RUN # 13
WORKFILE ID: C
WORKFILE NAME:

JAN/06/87 18:25:41

HEIGHT%

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.38	10938	BV	0.078	1.700
1.49	21932	D VV	0.107	3.418
1.66	31463	D VP	0.117	4.891
2.02	31501	PV	0.118	4.897
2.41	40251	VV	0.239	6.257
2.89	20815	VV	0.206	3.112
3.12	11907	VV	0.208	1.851
3.47	14083	VV	0.212	2.189
3.98	34550	VV	0.258	5.371
4.53	32138	VP	0.171	4.995
4.93	7962	PV	0.140	1.098
5.24	125135	VV	0.204	19.453
5.46	124083	VV	0.329	19.290
6.96	22377	VP	0.421	3.479
8.14	2396	PB	0.289	0.359
9.59	7168	BP	0.650	1.114
11.02	4212	PV	0.471	0.655
12.10	22296	VV	0.695	3.466
13.05	16632	VV	0.531	2.593
13.81	18600	VV	1.079	2.892
16.00	15309	VV	1.676	2.380
16.61	22791	VV	1.011	3.543
20.93	3026	VV	1.351	0.478
27.05	3390	PP	2.143	0.527

Page 1
Received: 12/09/86

SPECTRUM REPORT
Results by Sample

Work Order # 86-12-041

SAMPLE ID CAV-DWII-001 (LLW)		SAMPLE # 03 FRACTIONS: A		Date & Time Collected 12/09/86		Category					
AG	<50 ug/L	AS	522 ug/L	BE	<15 ug/L	CD	<25 ug/L	CN_SU	<5000 ug/L	CR	<45 ug/L
CU	<50 ug/L	FE	10500 ug/L	HG	<0.2 ug/L	NI	395 ug/L	N03_SU	N/A ug/L	PB	28.3 ug/L
SB	<220 ug/L	SE	<25 ug/L	TL	<50 ug/L	ZN	490 ug/L				

CHECKLIST

SECTION II - Sample Data Packets

Sample No: SCK-MW17-001

Organic Analysis Data Sheets (OADS)

VOA
 SV
 Pest.

Tentatively Identified Compounds

VOA
 SV

Raw Data

VOA chromatogram
 VOA quantitation report (surrogates)
 SV chromatogram
 SV quantitation report (surrogates)
 Pest. chromatogram and integrator print out

Inorganic Analysis Data Sheets (IADS)

003383

SPECTRIX DC # ---- 8

INST ID 4000

SAMPLE NUMBER: SCK-MW17-001

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

Laboratory Name: SPECTRIX
 Sample ID No.: 861204104
 Sample Matrix: WATER
 Data Release Authorized By: *[Signature]*

Case No.: ----
 QC Report No.: ----
 Contract No.: ----
 Date Sample Received: 12/9/86

VOLATILES

Concentration: Low
 Date Extracted/Prepared: 12/13/86
 Date Analyzed: 12/13/86

Datafile: 4U12041V04
 Dilution Factor: 1.0

CAS # COMPOUND

UG/L

003384

C010	CHLOROMETHANE	10 U
C015	BROMOMETHANE	10 U
C020	VINYL CHLORIDE	10 U
C025	CHLOROETHANE	10 U
C030	METHYLENE CHLORIDE	5 U
C035	ACETONE	10 U
C040	CARBON DISULFIDE	5 U
C045	1, 1-DICHLOROETHENE	5 U
C050	1, 1-DICHLOROETHANE	5 U
C055	TRANS-1, 2-DICHLOROETHENE	5 U
C060	CHLOROFORM	5 U
C065	1, 2-DICHLOROETHANE	5 U
C110	2-BUTANONE	10 U
C115	1, 1, 1-TRICHLOROETHANE	5 U
C120	CARBON TETRACHLORIDE	5 U
C125	VINYL ACETATE	10 U
C130	BROMODICHLOROMETHANE	5 U
C140	1, 2-DICHLOROPROPANE	5 U
C145	TRANS-1, 3-DICHLOROPROPENE	5 U
C150	TRICHLOROETHENE	5 U
C155	DIBROMOCHLOROMETHANE	5 U
C160	1, 1, 2-TRICHLOROETHANE	5 U
C165	BENZENE	5 U
C170	CIS-1, 3-DICHLOROPROPENE	5 U
C175	2-CHLOROETHYL VINYL ETHER	10 U
C180	BROMOFORM	5 U
C210	2-HEXANONE	10 U
C205	4-METHYL-2-PENTANONE	10 U
C220	TETRACHLOROETHENE	5 U
C225	1, 1, 2-TETRACHLOROETHANE	5 U
C230	TOLUENE	5 U
C235	CHLOROBENZENE	5 U
C240	ETHYLBENZENE	5 U
C245	STYRENE	5 U
C250	TOTAL XYLENES	5 U

= UNDETECTED AT THE LISTED DETECTION LIMIT

= COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

= COMPOUND ALSO FOUND IN BLANK

NST ID: 4530

SPECTRIX DC # ---- 8
SAMPLE NUMBER: SCK-MW17-001

ORGANICS ANALYSIS DATA SHEET - PAGE 2

LABORATORY NAME: SPECTRIX
 LAB SAMPLE ID NO.: 861204104
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *YPH*

CASE NO.: ---
 QC REPORT NO.: ---
 CONTRACT NO.: ---
 DATE SAMPLE RECEIVED: *12/9/86*

SEMICOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED: *12/13/86*
 DATE ANALYZED: 01/09/87

DATAFILE: 4U12041C04

DILUTION FACTOR:2....

CAS #	COMPOUND	UG/L	
C315	PHENOL	20 U	
C325	BIS(2-CHLOROETHYL)ETHER	20 U	5
C330	2-CHLOROPHENOL	20 U	80
C335	1, 3-DICHLOROBENZENE	20 U	5
C340	1, 4-DICHLOROBENZENE	20 U	5
C345	BENZYL ALCOHOL	20 U	3
C350	1, 2-DICHLOROBENZENE	20 U	3
C355	2-METHYLPHENOL	20 U	0
C360	BIS(2-CHLOROISOPROPYL)ETHER	20 U	
C365	4-METHYLPHENOL	20 U	
C370	N-NITROSODIPropylAMINE	20 U	
C375	HEXAChLORoETHANE	20 U	
C410	NITROBENZENE	20 U	
C415	ISOPHORONE	20 U	
C420	2-NITROPHENOL	20 U	
C425	2, 4-DIMETHYLPHENOL	20 U	
C430	BENZOIC ACID	100 U	
C435	BIS(2-CHLOROETHOXY)METHANE	20 U	
C440	2, 4-DICHLOROPHENOL	20 U	
C445	1, 2, 4-TRICHLOROBENZENE	20 U	
C450	NAPHTHALENE	1 U	
C455	4-CHLORoANILINE	20 U	
C460	HEXAChLORoBUTADIENE	20 U	
C465	P-CHLORo-M-CRESOL	20 U	
C470	2-METHYLNAPHTHALENE	20 U	
C510	HEXAChLORoCYCLOPENTADIENE	20 U	
C515	2, 4, 6-TRICHLOROPHENOL	20 U	
C520	2, 4, 5-TRICHLOROPHENOL	100 U	
C525	2-CHLORoNAPHTHALENE	20 U	
C530	2-NITROANILINE	100 U	
C535	DIMETHYL PHTHALATE	20 U	
C540	ACENAPHTHYLENE	20 U	
C545	3-NITROANILINE	100 U	
C550	ACENAPHTHENE	20 U	
C555	2, 4-DINITROPHENOL	100 U	
C560	4-NITROPHENOL	100 U	
C565	DIBENZOFURAN	20 U	
C570	2, 4-DINITROTOLUENE	20 U	
C575	2, 6-DINITROTOLUENE	20 U	

SPECTRIX DC # ---- 8

SAMPLE NUMBER: SCK-MW17-001

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

DATAFILE: 4U12041C04

CAS #	COMPOUND	UG/L	
C580	DIETHYL PHTHALATE	20	U
C585	4-CHLOROPHENYL PHENYL ETHER	20	U
C590	FLUORENE	20	U
C595	4-NITROANILINE	100	U
C610	4, 6-DINITRO-2-METHYLPHENOL	100	U
C615	N-NITROSODIPHENYLAMINE	20	U
C625	4-BROMOPHENYL PHENYL ETHER	20	U
C630	HEXAChLOROBENZENE	20	U
C635	PENTACHLOROPHENOL	100	U
C640	PHENANTHRENE	20	U
C645	ANTHRACENE	20	U
C650	DI-N-BUTYL PHTHALATE	20	U
C655	FLUORANTHENE	20	U
C715	PYRENE	20	U
C720	BUTYL BENZYL PHTHALATE	20	U
C725	3, 3'-DICHLOROBENZIDINE	40	U
C730	BENZO(A)ANTHRACENE	20	U
C735	BIS(2-ETHYLHEXYL)PHTHALATE	20	U
C740	CHRYSENE	20	U
C760	DI-N-OCTYL PHTHALATE	20	U
C765	BENZO(B)FLUORANTHENE	20	U
C770	BENZO(K)FLUORANTHENE	20	U
C775	BENZO(A)PYRENE	20	U
C780	INDENO(1, 2, 3-CD)PYRENE	20	U
C785	DIBENZO(A, H)ANTHRACENE	20	U
C790	BENZO(GHI)PERYLENE	20	U

003386

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

Lab Name: Spectrix
Case No: KEYSTONE
Lab No.: 86-12-041-04A

DC No: KEYSTONE
SMO No: SCK-MW17-001

ORGANICS ANALYSIS DATA SHEET
(page 3)

Pesticide/PCBs

Conc: Low GPC Cleanup Yes No
Date Extracted: 12/11/86 Sep. Funnel Ext. Yes
Date Analyzed: 12/23/86 Cont. Liquid Ext. Yes

Dilution Factor: 1

Percent Moisture(decanted):

CAS #	Compound	ug/L
319-84-6	ALPHA-BHC	0.05 u
319-85-7	BETA-BHC	0.05 u
319-86-8	DELTA-BHC	0.05 u
58-89-9	GAMMA-BHC (LINDANE)	0.05 u
76-44-8	HEPTACHLOR	0.05 u
309-00-2	ALDRIN	0.05 u
1024-57-3	HEPTACHLOR EPOXIDE	0.05 u
959-98-8	ENDOSULFAN I	0.05 u
60-57-1	DIELDRIN	0.10 u
72-55-9	4,4'-DDE	0.10 u
72-20-8	ENDRIN	0.10 u
33213-65-9	ENDOSULFAN II	0.10 u
72-54-8	4,4'-DDD	0.10 u
7421-93-4	ENDRIN ALDEHYDE	0.10 u
1031-07-8	ENDOSULFAN SULFATE	0.10 u
50-29-3	4,4'-DDT	0.10 u
72-43-5	METHOXYSCHLOR	0.50 u
53494-70-5	ENDRIN KETONE	0.10 u
57-74-9	CHLORDANE	0.50 u
8001-35-2	TOXAPHENE	1.0 u
12674-11-2	AROCLOL-1016	0.50 u
11104-28-2	AROCLOL-1221	0.50 u
11141-16-5	AROCLOL-1232	0.50 u
53469-21-9	AROCLOL-1242	0.50 u
12672-29-6	AROCLOL-1248	0.50 u
11097-69-1	AROCLOL-1254	1.0 u
11096-82-5	AROCLOL-1260	1.0 u

0037
0035
0035

V_s = 500 = Volume of water extracted (ml)

V_t = 5000 = Volume of total extract (uL)

V_i = 1 = Volume of extract injected(uL)

1st Column Inst. : 6000 N.A.

2nd Column Inst. : 6000

U=UNDETECTED AT THE LISTED DETECTION LIMIT.
J=COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT.
B=COMPOUND ALSO FOUND IN BLANK.

INST ID: 4000

SPECTRIX DC # ----- B

SAMPLE NUMBER: SCK-MW17-001

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: -----

QC REPORT NO.: -----

ANALYST: CKT

DATAFILE: 4U12041V04

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	VOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT	OO
124-38-9	CARBON DIOXIDE (ACN)*	34	968	28	100
124-38-9	CARBON DIOXIDE (ACN)*	48	977	34	100

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

* ALSO FOUND IN BLANK

INST ID: 4530

SPECTRIX DC # ---- 8

SAMPLE: SCK-MW17-001

ORGANICS ANALYSIS DATA SHEET - PAGE 5

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: ---

GC REPORT NO.: ---

ANALYST: GU

DATAFILE: 4U12041C04

B. TENTATIVELY IDENTIFIED COMPOUNDS

FAS #	SEMOVOLATILE COMPOUND NAMES		SCAN#	PURITY	AMOUNT
10-12-3	2-HEXANONE, 5-METHYL-	*	333	929	24 M
540-97-6	CYCLOHEXASILOXANE, DODECAMETHYL-	*	890	867	15 M
540-97-6	CYCLOHEXASILOXANE, DODECAMETHYL-	*	895	869	89

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

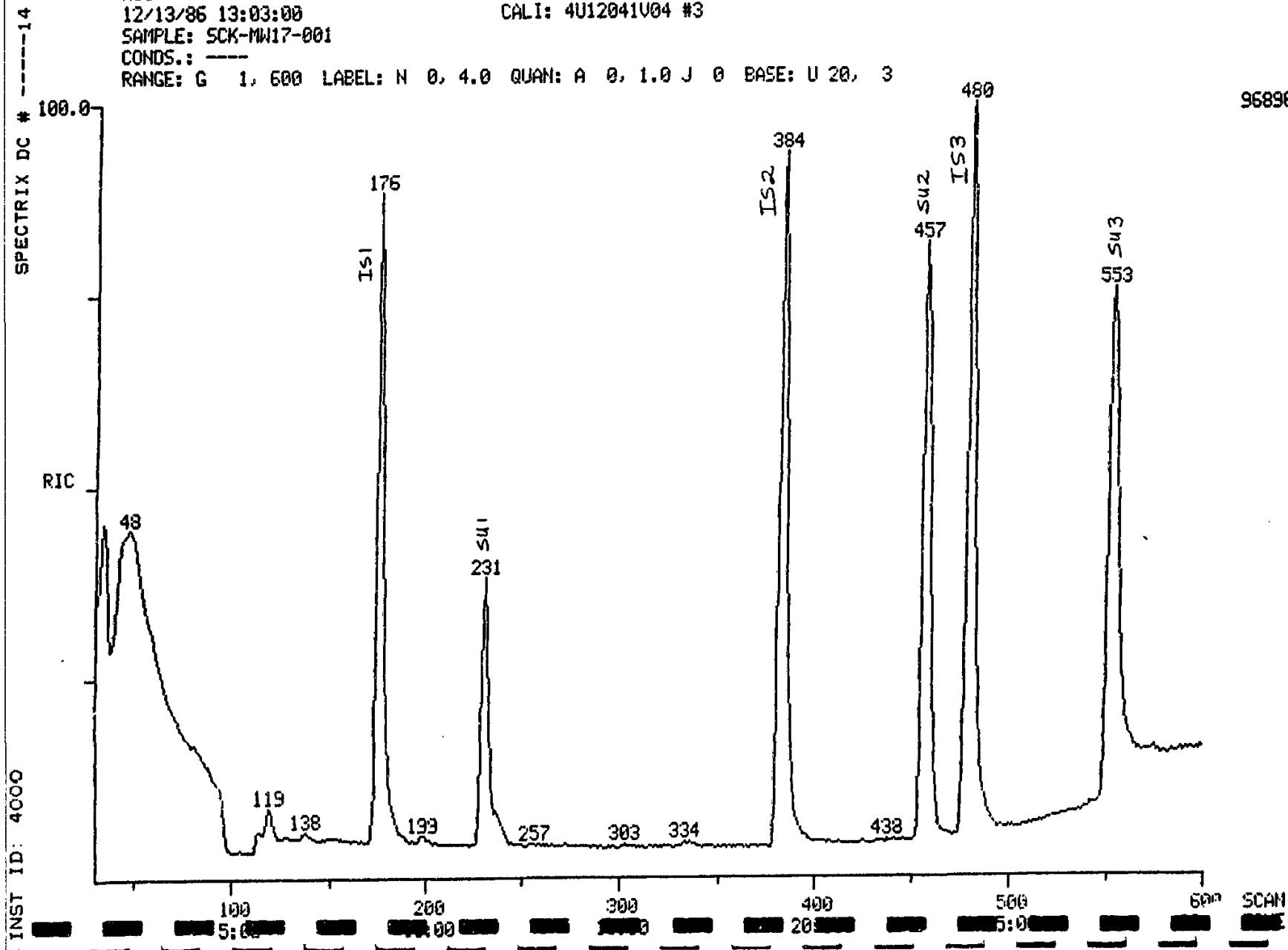
* ALSO FOUND IN BLANK

062200

RIC
12/13/86 13:03:00
SAMPLE: SCK-MW17-001
COND.:
RANGE: G 1, 600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12041U04 #1 SCANS 30 TO 600
CALI: 4U12041U04 #3

96896.



SPECTRIX DC # -----14

VDA QUANTITATION REPORT

SAMPLE ID.	SCK-MW17-001		
.ENAME	4U12041V04	INST	4000
CLIENT	KEYSTONE	ANALYST	CKT
DATE INJECTED	12/13/86 13:03:00	VERIFIED BY	<i>YTH</i>
ALIB STD DATE	12-14-86	CORR. FACTOR	1.00

CAN#	VDA COMPOUNDS	M/E	AMOUNT	RRT	AREA
176 CI01	BROMOCHLOROMETHANE	**IS1**	128	50 UG/L	1.000 34542
383 CI10	1,4-DIFLUOROBENZENE	**IS2**	114	50 UG/L	1.000 146666
479 CI20	D5-CHLOROBENZENE	**IS3**	117	50 UG/L	1.000 120310
231 CS15	D4-1,2-DICHLOROETHANE	**SU1**	65	95 %	1.313 33604
457 CS05	D8-TOLUENE	**SU2**	98	99 %	0.954 98643
553 CS10	P-BROMOFLUOROBENZENE	**SU3**	95	98 %	1.154 609510
119 CO30	METHYLENE CHLORIDE *		84	3 UG/L	0.676 2169M
238 C110	2-BUTANONE *		72	8 UG/L	0.621 1179M

* BLANK CORRECTED ON DATA SHEET

0
0

263390

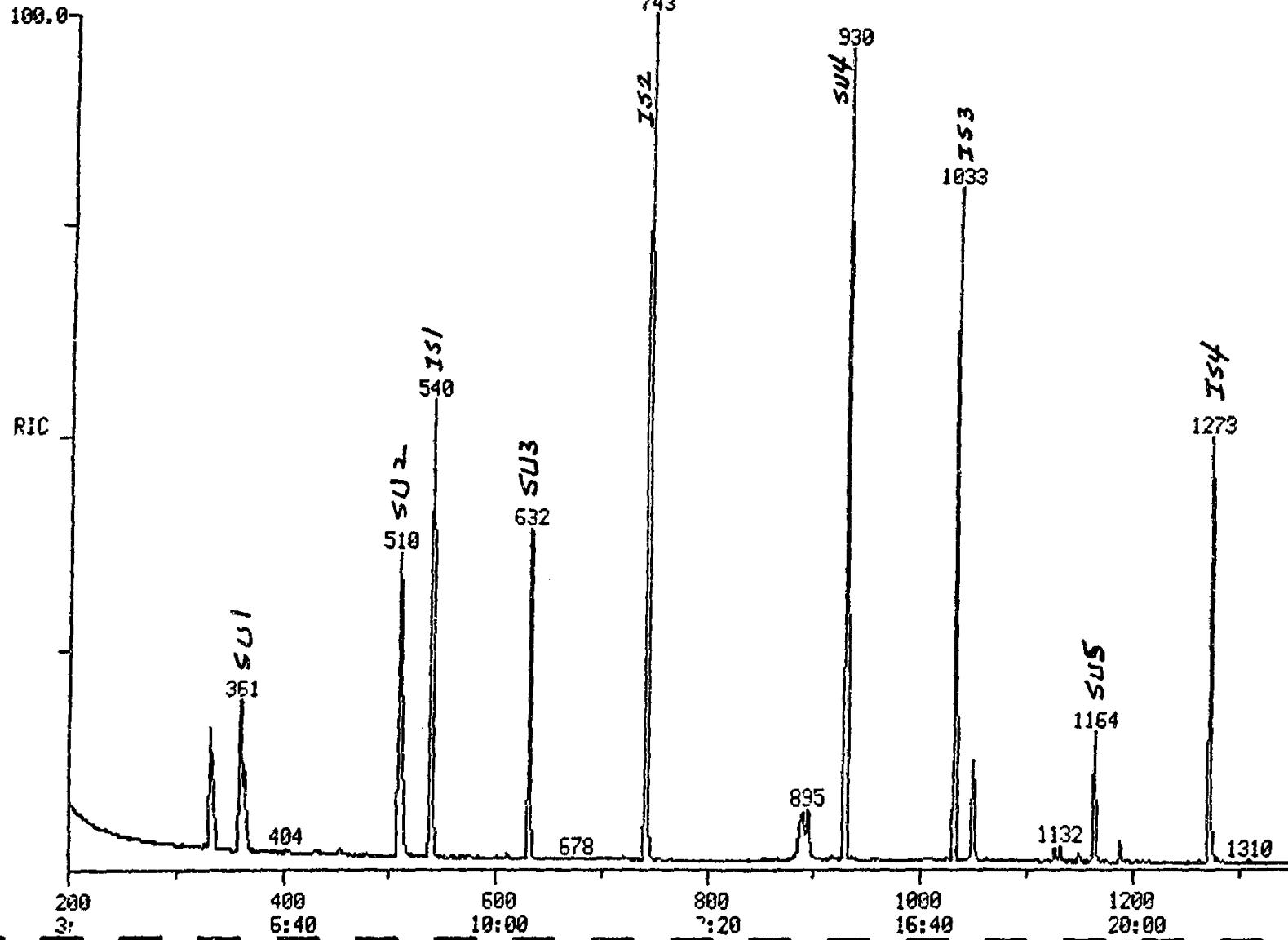
RIC
01/09/87 14:02:00
SAMPLE: SCK-MN17-001
COND.: ---
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12041C04 #1
CALI: 4U12041C04 #3

SCANS 200 TO 1350

SPECTRIX DC # -----14

INST ID: 4530



99456.

003392

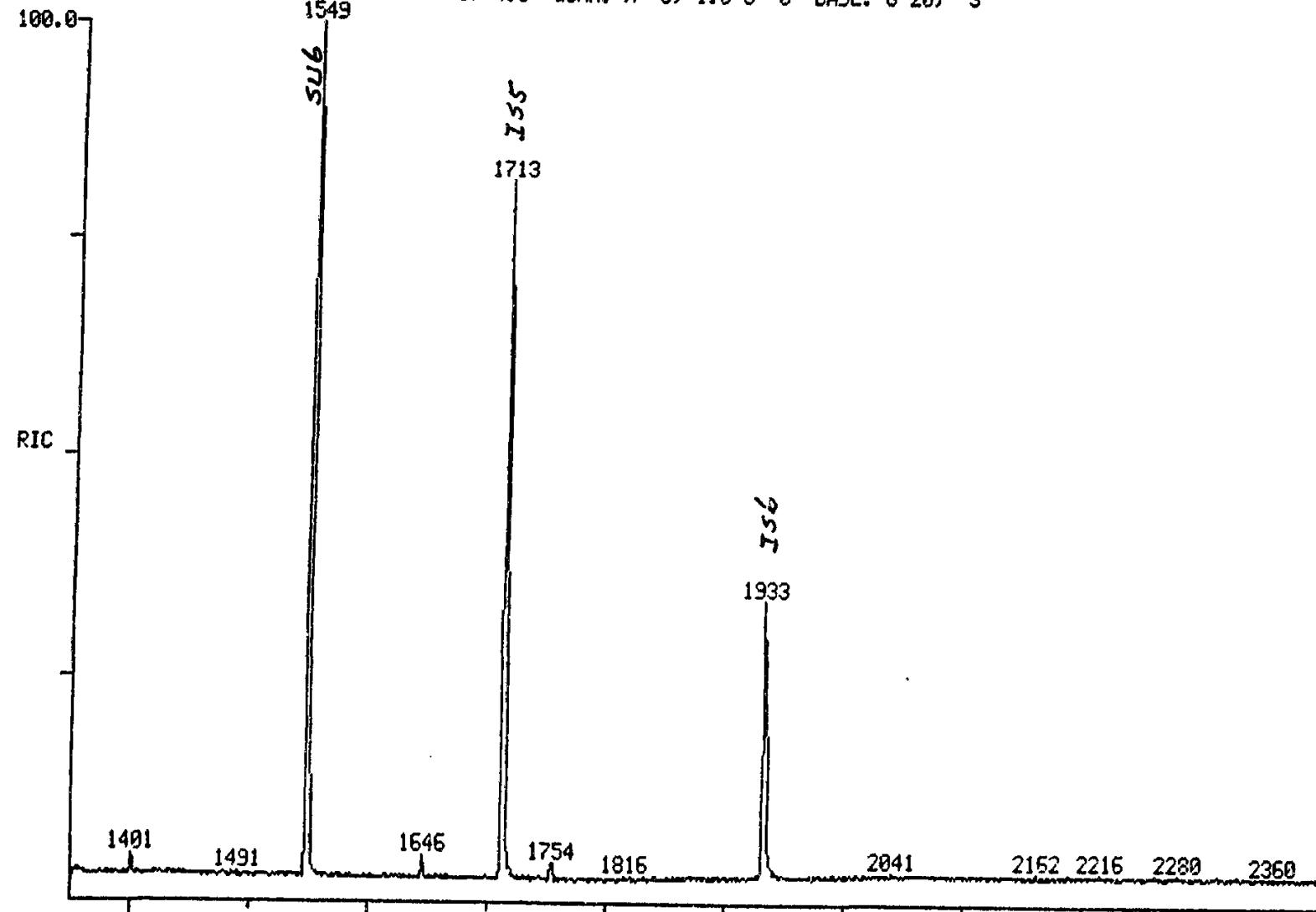
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RIC
01/09/87 14:02:00
SAMPLE: SCK-MW17-001
COND.: ---
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12041C04 #1 SCANS 1350 TO 2400
CALI: 4U12041C04 #3

25888.

SPECTRIX DC # ----14



INSTR ID: 4530

003393

SPECTRIX DC # ----14

SEMIVOLATILE WATER REPORT

SAMPLE ID. SCK-MW17-001

.LENAME 4U12041C04

INST ID. 4530

CLIENT KEYSTONE

ANALYST GU

DATE INJECTED 01/09/87 14:02:00

VERIFIED BY JPH

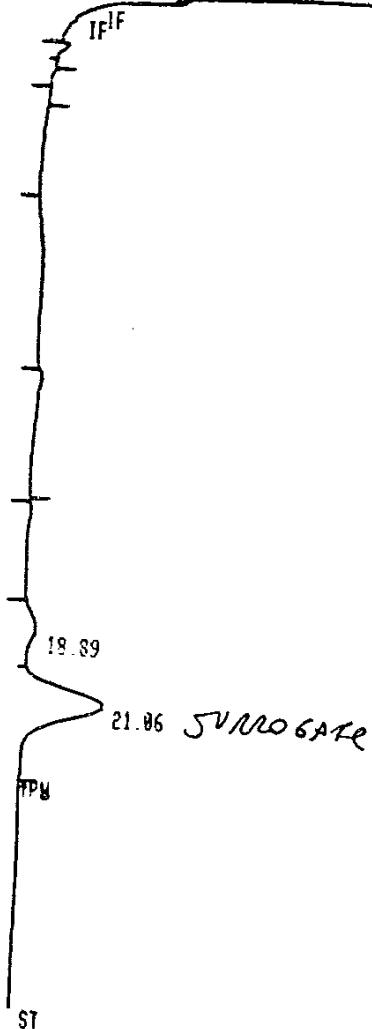
CALIB STD DATE 01/08/87

CORR. FACTOR 2.00

SCAN#	SEMIVOLATILE COMPOUNDS	M/E	AMOUNT	RRT	AREA
540	CI30 1, 4-DICHLOROBENZEND-D4 ** IS1 **	152	80 UG/L	1. 000	2425
743	CI40 NAPHTHALENE-D8 ** IS2 **	136	80 UG/L	1. 000	9617
1033	CI50 ACENAPHTHENE-D10 ** IS3 **	164	80 UG/L	1. 000	4771
1273	CI60 PHENANTHRENE-D10 ** IS4 **	188	80 UG/L	1. 000	5742
1713	CI70 CHRYSENE-D12 ** IS5 **	240	80 UG/L	1. 000	2760
1933	CI75 PERYLENE-D12 ** IS6 **	264	80 UG/L	1. 000	1482
360	CS50 2-FLUOROPHENOL ** SU1 **	112	58 %	0. 667	2362
510	CS45 PHENOL-D5 ** SU2 **	99	72 %	0. 944	4127
1164	CS55 2, 4, 6-TRIBROMOPHENOL ** SU5 **	330	53 %	1. 127	342
632	CS20 NITROBENZENE-D5 ** SU3 **	82	114 %	0. 851	308
930	CS25 2-FLUOROBIPHENYL ** SU4 **	172	115 %	0. 900	672
1549	CS30 TERPHENYL-D14 ** SU6 **	244	152 %	0. 904	2807
746	C450 NAPHTHALENE	128	1 UG/L	1. 004	169
1127	C580 DIETHYL PHTHALATE *	149	3 UG/L	1. 091	180

* BLANK CORRECTED ON DATA SHEET

STAMP: 86-12-Q41-Q4A ROC



DC #: N/A
SPECTRIX CORPORATION
CASE NO. Kayrone
SAMPLE NO. SCK-MW17-001
PESTICIDE FRACTION
1st Column - 1D
1.8m x 4mm ID x 6mm 00 5
1.5% SP 2250 5
1.95% SP2401 0 0 3
100/120 Supelcoport
Column Temp. 203 °C
Volume Inj. 1.0 ul
Instrument ID VARIAN 60

RUN #: 138
WORKFILE ID: C DEC/23/86 21:50:22
WORKFILE NAME:

HEIGHT%
RT HEIGHT TYPE AR/HT HEIGHT%
18.89 1553 BY 1.004 13.277
21.06 18144 VB 1.020 86.723

TOTAL HIGHT= 11697
MUL FACTOR= 1.0000E+00

Page 1

Received: 12/09/86

SPECTRIS 200 REPORT
Results by Sample

WORK ORDER # 86-12-041

SAMPLE ID <u>SCK-MW17-001 (LLW)</u>		SAMPLE # <u>04</u> FRACTIONS: A		Date & Time Collected <u>12/09/86</u>		Category _____	
AG	<10 ug/L	AS	<10 ug/L	BE	<3 ug/L	CD	<5 ug/L
CU	<10 ug/L	FE	<30 ug/L	HG	<0.2 ug/L	NI	38 ug/L
SB	<44 ug/L	SE	<5 ug/L	TL	<10 ug/L	ZN	27 ug/L

CHECKLIST

SECTION II - Sample Data Packets

Sample No: SCK-P03-001

Organic Analysis Data Sheets (OADS)

VOA
 SV
 Pest.

Tentatively Identified Compounds

VOA
 SV

Raw Data

VOA chromatogram
 VOA quantitation report (surrogates)
 SV chromatogram
 SV quantitation report (surrogates)
 Pest. chromatogram and integrator print out

Inorganic Analysis Data Sheets (IADS)

003397

SPECTRIX DC # ----- 8

INST ID 4000

SAMPLE NUMBER: SCK-P03-001

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

LABORATORY NAME: SPECTRIX
 SAMPLE ID NO.: 861204105
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *[Signature]*

CASE NO.: -----
 GC REPORT NO.: -----
 CONTRACT NO.: -----
 DATE SAMPLE RECEIVED: 12/13/86

VOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED/PREPARED: 12/13/86
 DATE ANALYZED: 12/13/86

DATAFILE: 4U12041V05
 DILUTION FACTOR: .1000.....

CAS # COMPOUND

UG/L

003398

C010	CHLOROMETHANE	1000 U
C015	BROMOMETHANE	1000 U
C020	VINYL CHLORIDE	1000 U
C025	CHLOROETHANE	1000 U
C030	METHYLENE CHLORIDE	500 U
C035	ACETONE	1000 J
C040	CARBON DISULFIDE	500 U
C045	1, 1-DICHLOROETHENE	500 U
C050	1, 1-DICHLOROETHANE	500 U
C055	TRANS-1, 2-DICHLOROETHENE	500 U
C060	CHLOROFORM	500 U
C065	1, 2-DICHLOROETHANE	500 U
C110	2-BUTANONE	1000 U
C115	1, 1, 1-TRICHLOROETHANE	500 U
C120	CARBON TETRACHLORIDE	500 U
C125	VINYL ACETATE	1000 U
C130	BROMODICHLOROMETHANE	500 U
C140	1, 2-DICHLOROPROPANE	500 U
C145	TRANS-1, 3-DICHLOROPROPENE	500 U
C150	TRICHLOROETHENE	500 U
C155	DIBROMOCHLOROMETHANE	500 U
C160	1, 1, 2-TRICHLOROETHANE	500 U
C165	BENZENE	800 U
C170	CIS-1, 3-DICHLOROPROPENE	500 U
C175	2-CHLOROETHYL VINYL ETHER	1000 U
C180	BROMOFORM	500 U
C210	2-HEXANONE	1000 U
C205	4-METHYL-2-PENTANONE	1000 U
C220	TETRACHLOROETHENE	500 U
C225	1, 1, 2-TETRACHLOROETHANE	500 U
C230	TOLUENE	740 U
C235	CHLOROBENZENE	500 U
C240	ETHYLBENZENE	500 U
C245	STYRENE	500 U
C250	TOTAL XYLEMES	500 J

- = UNDETECTED AT THE LISTED DETECTION LIMIT
- = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT
- = COMPOUND ALSO FOUND IN BLANK

INST ID: 4530

SPECTRIX DC # ---- 8
SAMPLE NUMBER: SCK-P03-001

ORGANICS ANALYSIS DATA SHEET - PAGE 2

BORATORY NAME: SPECTRIX
 LAB SAMPLE ID NO.: B61204105
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *[Signature]*

CASE NO.: ---
 QC REPORT NO.: ---
 CONTRACT NO.: ---
 DATE SAMPLE RECEIVED: 12/19/86

SEMOVOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED: 12/13/86
 DATE ANALYZED: 01/09/87

DATAFILE: 4U12041C05
 DILUTION FACTOR: 1.000

CAS #	COMPOUND	UG/L	
C315	PHENOL	10000 U	0
C325	BIS(2-CHLOROETHYL)ETHER	10000 U	6
C330	2-CHLOROPHENOL	10000 U	6
C335	1, 3-DICHLOROBENZENE	10000 U	5
C340	1, 4-DICHLOROBENZENE	10000 U	3
C345	BENZYL ALCOHOL	10000 U	3
C350	1, 2-DICHLOROBENZENE	10000 U	3
C355	2-METHYLPHENOL	10000 U	3
C360	BIS(2-CHLOROISOPROPYL)ETHER	10000 U	3
C365	4-METHYLPHENOL	10000 U	3
C370	N-NITROSODIPOXYLAMINE	10000 U	3
C375	HEXACHLOROETHANE	10000 U	3
C410	NITROBENZENE	10000 U	3
C415	ISOPHORONE	10000 U	3
C420	2-NITROPHENOL	10000 U	3
C425	2, 4-DIMETHYLPHENOL	10000 U	3
C430	BENZOIC ACID	50000 U	3
C435	BIS(2-CHLOROETHOXY)METHANE	10000 U	3
C440	2, 4-DICHLOROPHENOL	10000 U	3
C445	1, 2, 4-TRICHLOROBENZENE	10000 U	3
C450	NAPHTHALENE	140000	3
C455	4-CHLOROANILINE	10000 U	3
C460	HEXACHLOROBUTADIENE	10000 U	3
C465	P-CHLORO-M-CRESOL	10000 U	3
C470	2-METHYLNAPHTHALENE	19000	3
C510	HEXACHLOROCYCLOPENTADIENE	10000 U	3
C515	2, 4, 6-TRICHLOROPHENOL	10000 U	3
C520	2, 4, 5-TRICHLOROPHENOL	50000 U	3
C525	2-CHLORONAPHTHALENE	10000 U	3
C530	2-NITROANILINE	50000 U	3
C535	DIMETHYL PHTHALATE	10000 U	3
C540	ACENAPHTHYLENE	10000 U	3
C545	3-NITROANILINE	50000 U	3
C550	ACENAPHTHENE	26000	3
C555	2, 4-DINITROPHENOL	50000 U	3
C560	4-NITROPHENOL	50000 U	3
C565	DIBENZOFURAN	18000	3
C570	2, 4-DINITROTOLUENE	10000 U	3
C575	2, 6-DINITROTOLUENE	10000 U	3

SPECTRIX DC # ---- 8

SAMPLE NUMBER: SCK-P03-001

EMIVOLATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

DATAFILE: 4U12041C05

CAS #	COMPOUND	UG/L
C580	DIETHYL PHTHALATE	10000 U
C585	4-CHLOROPHENYL PHENYL ETHER	10000 U
C590	FLUORENE	19000
C595	4-NITROANILINE	50000 U
C610	4, 6-DINITRO-2-METHYLPHENOL	50000 U
C615	N-NITROSODIPHENYLAMINE	10000 U
C625	4-BROMOPHENYL PHENYL ETHER	10000 U
C630	HEXACHLOROBENZENE	10000 U
C635	PENTACHLOROPHENOL	50000 U
C640	PHENANTHRENE	58000
C645	ANTHRACENE	8300 J
C650	DI-N-BUTYL PHTHALATE	10000 U
C655	FLUORANTHENE	26000
C715	PYRENE	19000
C720	BUTYL BENZYL PHTHALATE	10000 U
C725	3, 3'-DICHLOROBENZIDINE	20000 U
C730	BENZO(A)ANTHRACENE	5000 J
C735	BIS(2-ETHYLHEXYL)PHTHALATE	10000 U
C740	CHRYSENE	4900 J
C760	DI-N-OCTYL PHTHALATE	10000 U
C765	BENZO(B)FLUORANTHENE	2200 J
C770	BENZO(K)FLUORANTHENE	2600 J
C775	BENZO(A)PYRENE	2600 J
C780	INDENO(1, 2, 3-CD)PYRENE	10000 U
C785	DIBENZO(A, H)ANTHRACENE	10000 U
C790	BENZO(GHI)PERYLENE	10000 U

003400

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

Lab Name: Spectrix
Case No: KEYSTONE
Lab No.: 86-12-041-05A

DC No: KEYSTONE
SMO No: SCK-P03-001

ORGANICS ANALYSIS DATA SHEET
(page 3)

Pesticide/PCBs

Conc: Low
Date Extracted: 12/11/86
Date Analyzed: 12/23/86
Dilution Factor: 1
Percent Moisture(decant):

GPC Cleanup Yes No
Sep. Funnel Ext. Yes
Cont. Liquid Ext. Yes

CAS #	Compound	ug/L
319-84-6	ALPHA-BHC	0.05 U
319-85-7	BETA-BHC	0.05 U
319-86-8	DELTA-BHC	0.05 U
58-89-9	GAMMA-BHC (LINDANE)	0.05 U
76-44-8	HEPTACHLOR	0.05 U
309-00-2	ALDRIN	0.05 U
1024-57-3	HEPTACHLOR EPOXIDE	0.05 U
959-98-8	ENDOSULFAN I	0.05 U
60-57-1	DIELDRIN	0.10 U
72-55-9	4,4'-DDE	0.10 U
72-20-8	ENDRIN	0.10 U
33213-65-9	ENDOSULFAN II	0.10 U
72-54-8	4,4'-DDD	0.10 U
7421-93-4	ENDRIN ALDEHYDE	0.10 U
1031-07-8	ENDOSULFAN SULFATE	0.10 U
50-29-3	4,4'-DDT	0.10 U
72-43-5	METHOXYPHENYLCHLOR	0.50 U
53494-70-5	ENDRIN KETONE	0.10 U
57-74-9	CHLORDANE	0.50 U
8001-35-2	TOXAPHENE	1.0 U
12674-11-2	AROCLOTR-1016	0.50 U
11104-28-2	AROCLOTR-1221	0.50 U
11141-16-5	AROCLOTR-1232	0.50 U
53469-21-9	AROCLOTR-1242	0.50 U
12672-29-6	AROCLOTR-1248	0.50 U
11097-69-1	AROCLOTR-1254	1.0 U
11096-82-5	AROCLOTR-1260	1.0 U

003401

V_s = 500 = Volume of water extracted (ml)

V_t = 5000 = Volume of total extract (ul)

V_i = 1 = Volume of extract injected(ul)

1st Column Inst. : 5000
2nd Column Inst. : 6000 NA.

U=UNDETECTED AT THE LISTED DETECTION LIMIT.

J=COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT.

B=COMPOUND ALSO FOUND IN BLANK.

Form 1

INST ID: 4000

SPECTRIX DC # ----- 8

SAMPLE NUMBER: SCK-P03-001

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: -----

GC REPORT NO.: -----

ANALYST: TTO

DATAFILE: 4U12041V05

B. TENTATIVELY IDENTIFIED COMPOUNDS

AS #	VOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
124-38-9	CARBON DIOXIDE (ACN) *	49	970	3200 UG/L 0.0354

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

* ALSO FOUND IN BLANK

SAMPLE: SCK-P03-001

ORGANICS ANALYSIS DATA SHEET - PAGE 5

LABORATORY NAME: SPECTRIX CORPORATION CASE NO.: ---

GC REPORT NO.: --- ANALYST: GU DATAFILE: 4U12041C05

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	SEMIVOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
611-15-4	BENZENE, 1-ETHENYL-2-METHYL-	560	962	5700~
95-13-6	1H-INDENE	571	946	7800~
95-15-8	BENZO[B]THIOPHENE	749	925	6900~
2471-83-2	1H-INDENE, 1-ETHYLLIDENE-	874	917	9600~
92-52-4	1, 1'-BIPHENYL	939	932	5900 J
-81-40-8	NAPHTHALENE, 2, 3-DIMETHYL-	977	936	5500 J
132-65-0	DIBENZOTHIOPHENE	1254	770	4700 J
86-74-8	9H-CARBAZOLE	1314	897	4500 J
203-64-5	4H-CYCLOPENTA[DEF]PHENANTHRENE	1374	861	5800 J

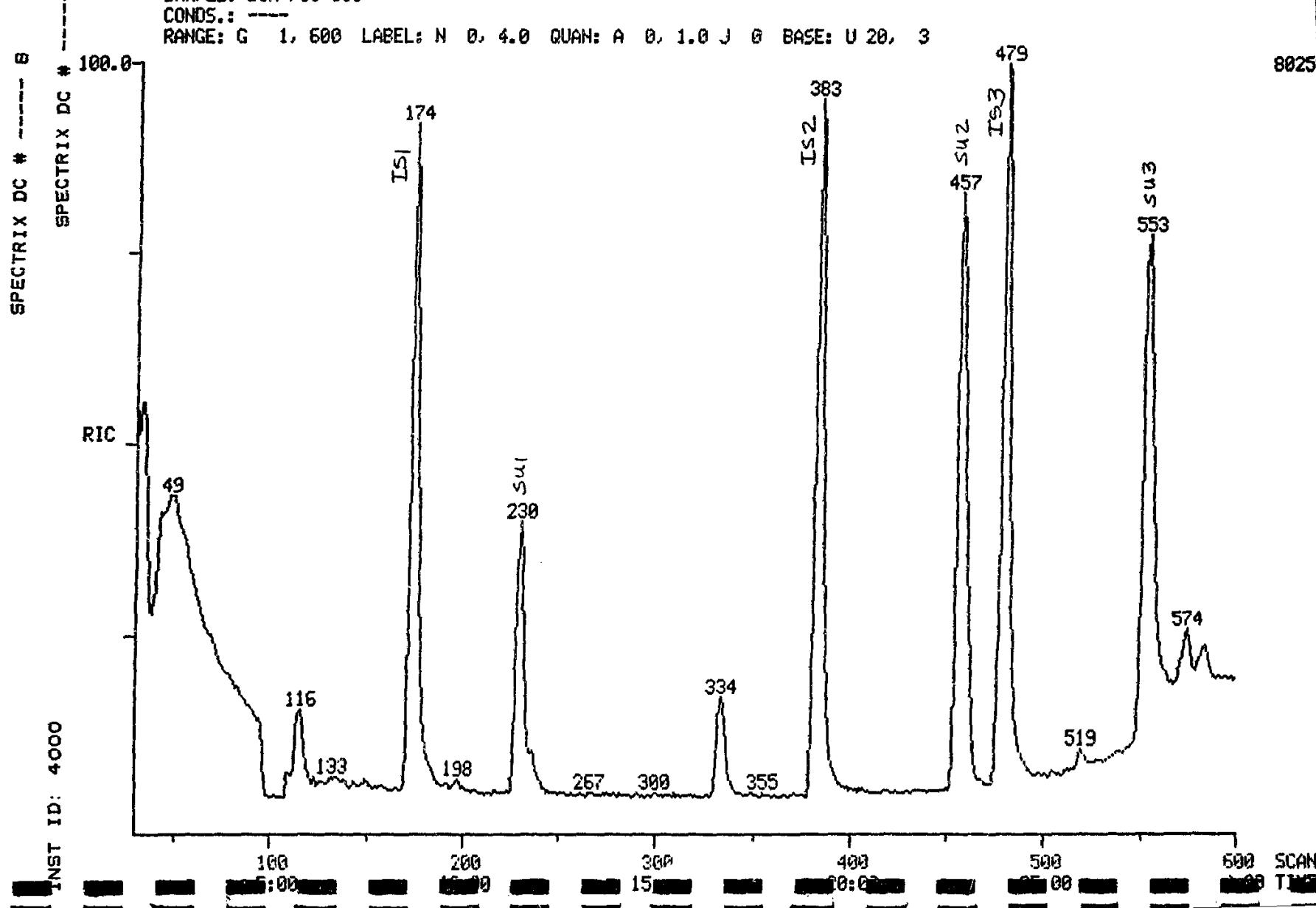
J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

003404

RIC
12/13/86 16:15:00
SAMPLE: SCK-P03-001
COND.: ---
RANGE: G 1, 600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12041V05 #1
CALI: 4U12041V05 #3

SCANS 30 TO 600



003404

SPECTRIX DC # -----14

VOA QUANTITATION REPORT

SAMPLE ID.	SCK-P03-001		
FILENAME	4U12041V05	INST	4000
CLIENT	KEYSTONE	ANALYST	TTO
DATE INJECTED	12/13/86 16:15:00	VERIFIED BY	<i>ypk</i>
ALIB STD DATE	121486	CORR. FACTOR	100.00

CAN#	VOA COMPOUNDS	M/E	AMOUNT	RRT	AREA
174 C101	BROMOCHLOROMETHANE	**IS1**	128	5000 UG/L	1.000 29882
383 C110	1, 4-DIFLUOROBENZENE	**IS2**	114	5000 UG/L	1.000 118884
479 C120	D5-CHLOROBENZENE	**IS3**	117	5000 UG/L	1.000 98380
230 CS15	D4-1, 2-DICHLOROETHANE	**SU1**	65	96 %	1.322 29553
457 CS05	D8-TOLUENE	**SU2**	98	98 %	0.954 79547
553 CS10	P-BROMOFLUOROBENZENE	**SU3**	95	96 %	1.154 49112
116 C030	METHYLENE CHLORIDE *		84	658 UG/L	0.667 4506
134 C035	ACETONE		43	416 UG/L	0.770 1039
237 C110	2-BUTANONE*		72	1105 UG/L	0.619 1273
333 C165	BENZENE		78	798 UG/L	0.869 17187
460 C230	TOLUENE		92	736 UG/L	0.960 9454
583 C250	TOTAL XYLEMES		106	347 UG/L	1.217 2861

* BLANK CORRECTED ON DATA SHEET

003406

RIC
01/09/87 18:41:00
SAMPLE: SCK-P03-001
COND.: ---
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12041C05 #1 SCANS 200 TO 1350
CALI: 4U12041C05 #3

265216.

SPECTRIX DC # -----14

100.0

RIC

INST ID: 4530

200

3:

400
6:40

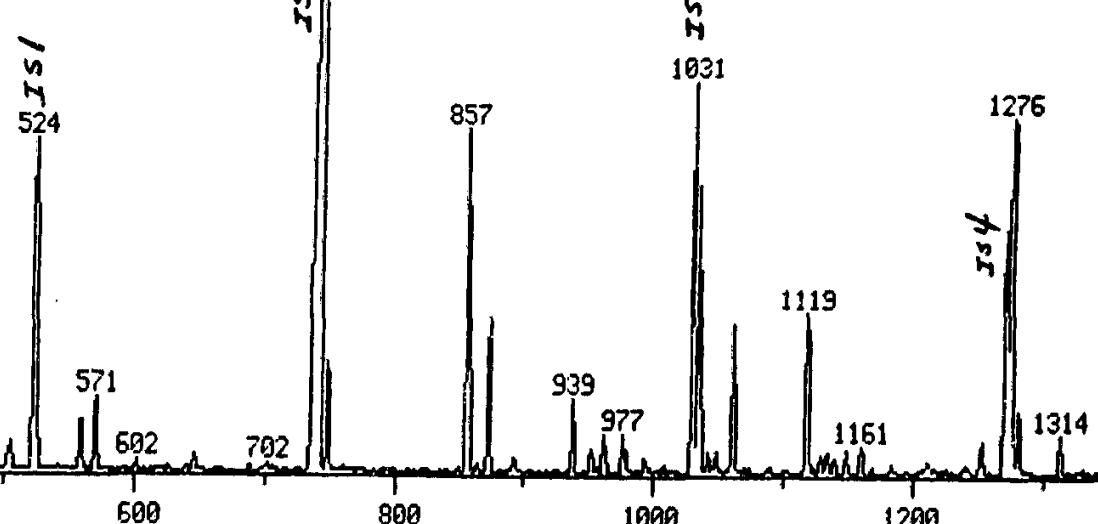
600
10:09

800
12:20

1000
16:40

1200
20:00

SCAN
TIME



003406

L O G E O

RIC
01/09/67 18:41:00
SAMPLE: SCK-P03-001
COND.: ---
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAH: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12041C05 #1 SCANS 1350 TO 2400
CALI: 4U12041C05 #3

35264.

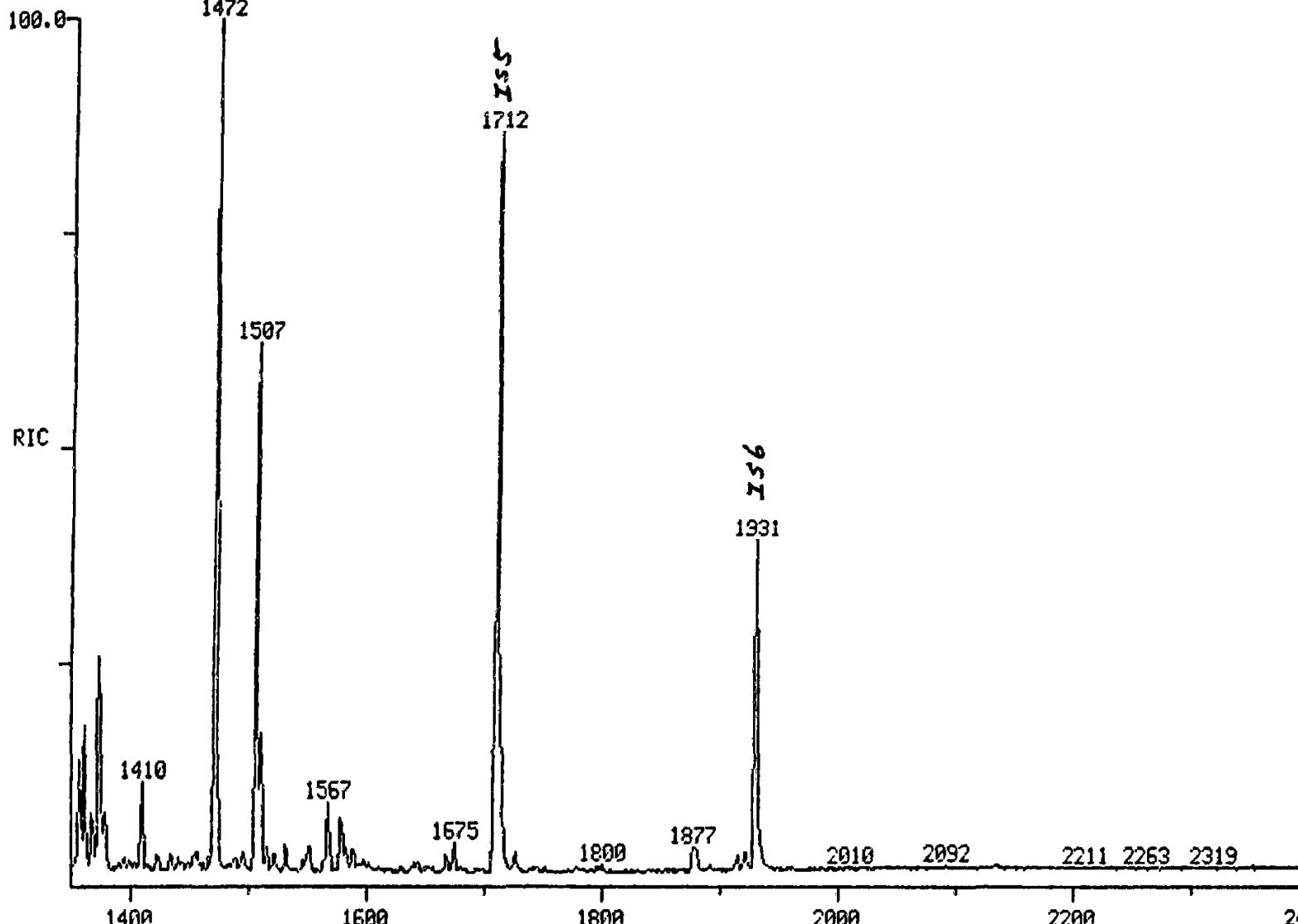
SPECTRIX DC # -----14

100.0

RIC

INST ID: 4530

003407



2400 SCAN
00 TIME

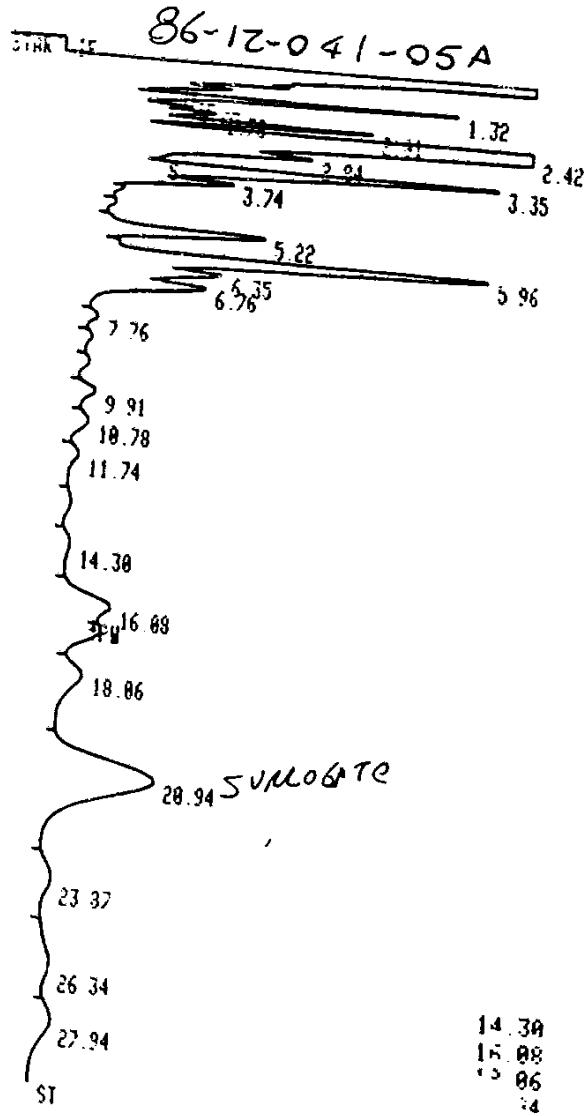
SPECTRIX DC # ----14

SEMIVOLATILE WATER REPORT

SAMPLE ID.	SCK-P03-001		
FILENAME	4U12041C05	INST ID.	4530
CLIENT	KEYSTONE	ANALYST	GU
DATE INJECTED	01/09/87 18:41:00	VERIFIED BY	<i>YPA</i>
CALIB STD DATE	01/08/87	CORR. FACTOR	1000.00

SCAN#	SEMIVOLATILE COMPOUNDS	M/E	AMOUNT	RRT	AREA
524	CI30 1, 4-DICHLOROBENZEND-D4 ** IS1 **	152	40000 UG/L	1.000	3511
738	CI40 NAPHTHALENE-D8 ** IS2 **	136	40000 UG/L	1.000	13828
1031	CI50 ACENAPHTHENE-D10 ** IS3 **	164	40000 UG/L	1.000	6784
1272	CI60 PHENANTHRENE-D10 ** IS4 **	188	40000 UG/L	1.000	8604
1712	CI70 CHRYSENE-D12 ** IS5 **	240	40000 UG/L	1.000	4208
1931	CI75 PERYLENE-D12 ** IS6 **	264	40000 UG/L	1.000	2328
743	C430 NAPHTHALENE	128	142377 UG/L	1.007	47273
857	C470 2-METHYLNAPHTHALENE	142	19483 UG/L	1.161	6006
1036	C550 ACENAPHTHENE	153	25876 UG/L	1.005	4033
1063	C565 DIBENZOFURAN	168	18030 UG/L	1.031	3812
1119	C590 FLUORENE	166	18545 UG/L	1.085	3282
1276	C640 PHENANTHRENE	178	58297 UG/L	1.003	12075
1282	C645 ANTHRACENE	178	8315 UG/L	1.008	1749
1472	C655 FLUORANTHENE	202	25557 UG/L	1.157	4950
1507	C715 PYRENE	202	19047 UG/L	0.880	3078
1709	C730 BENZO(A)ANTHRACENE	228	5027 UG/L	0.998	550
1715	C740 CHRYSENE	228	4905 UG/L	1.002	500
1877	C765 BENZO(B)FLUORANTHENE	252	2220 UG/L	0.972	152
1879	C770 BENZO(K)FLUORANTHENE	252	2623 UG/L	0.973	219
1921	C775 BENZO(A)PYRENE	252	2604 UG/L	0.995	157

* SURROGATES WERE DILUTED OUT



RUN # 139
WORKFILE ID: C
WORKFILE NAME:

DEC/23/86 22:21:38

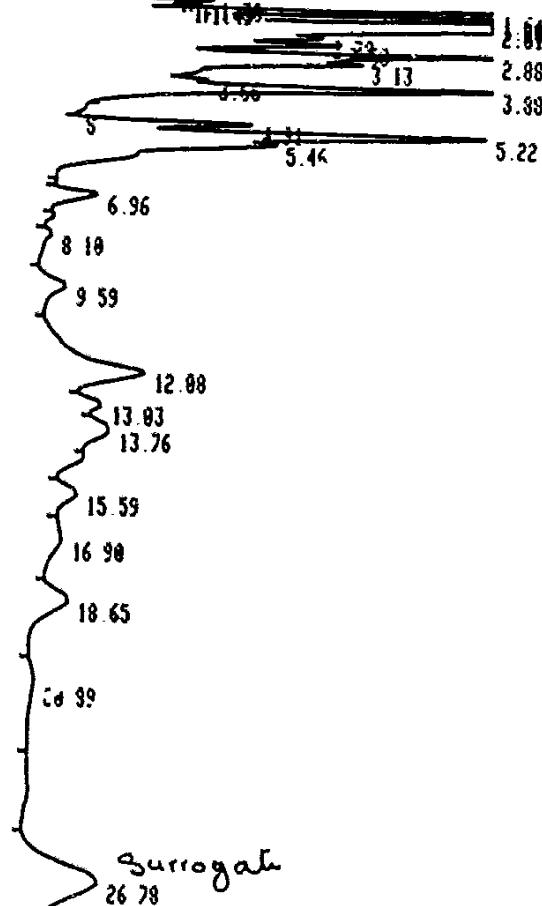
DC # N/A
SPECTRUM CORPORATION
CASE NO. Keystone
SAMPLE NO. SCK - P03-001
PESTICIDE FRACTION
1st Column - 1D
1.8m x 4mm ID x 6mm 00 0
1.5% SP 2250 0
1.95% SP2401 0
100/120 Supelcoport 0
Column Temp. 203 °C
Volume Inj. 1.0 ul
Instrument ID VARIAN 60

14.38	1155	PV	0.292	0.040
15.08	6714	VV	0.290	0.232
15.86	3782	VV	1.025	0.131
16.88	14017	VP	1.264	0.484
18.86	1672	PV	1.003	0.058
20.94	2124	VV	1.478	0.873
23.37	2728	I VP	1.393	0.094

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.32	49341	D BV	0.108	1.392
1.63	5718	VV	0.103	0.197
1.78	8893	VV	0.159	0.307
2.01	29654	D VH	0.108	1.023
2.42	2603560	SHB	0.141	89.827
2.84	10676	DTBB	0.125	0.368
3.35	45346	BV	0.210	1.565
3.74	13239	VP	0.162	0.457
5.22	20962	PV	0.247	0.723
5.96	50446	VV	0.271	1.741
6.35	16167	VV	0.289	0.558
6.76	14425	VP	0.337	0.498
7.76	1270	PV	0.298	0.044
9.91	2427	PV	0.447	0.084
10.78	1945	VP	0.558	0.067
11.74	1155	PV	0.421	0.049

STAX-# 86-12-041-06A

NA.



DC# _____
SPECTRIX HOUSTON

CASE NO. Keystone

SAMPLE NO. Sck-P03-00
PESTICIDE FRACTION
2ND COLUMN CONFIRMATI

2ND COLUMN TIC
1.8m x2mm ID x6mm OD
3% SP-2100
100/120 SUPELCOPORT
COLUMN TEMP. 187 °C
VOLUME INJ. 1 uL
INSTRUMENT ID. 6000

000
000

RUN # 14 JAN/06/87 19:01:57
WORKFILE ID: C
WORKFILE NAME:

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.38	23981	BV	0.093	0.629
1.49	28462	D VV	0.088	0.537
1.64	132795	D VV	0.092	3.485
1.88	198370	D VH	0.093	5.206
2.01	2813274	SHB	0.151	23.828
2.38	9556	DTBP	0.099	0.251
2.58	25378	DTPP	0.111	0.666
2.88	78415	TPV	0.177	2.058
3.13	47551	DTVV	0.168	1.248
3.66	17583	TVV	0.115	0.461
3.88	99802	TVB	0.207	2.619
4.91	45751	BV	0.188	1.201
5.22	102817	VV	0.224	2.829
5.46	53640	D VP	0.345	1.403
5.96	11657	VV	0.324	0.306
8.10	2144	PP	0.407	0.856
9.59	7190	PV	0.644	0.189
12.08	27576	VV	0.948	0.724
13.03	16937	VV	0.563	0.445
13.76	19127	VV	0.874	0.502
15.59	11428	VV	0.828	0.301
16.98	8054	VV	1.417	0.211
18.65	10275	VV	0.939	0.270
20.89	2109	VV	1.527	0.055
26.78	19679	PB	1.563	0.516

Page 1
Received: 12/09/86

SPECTRI[®] S O C REPORT
Results by Sample

Work Order # 86-12-041

SAMPLE ID <u>SCK-P03-001 (LLW)</u>		SAMPLE # <u>05</u>		FRACTIONS: <u>A</u>			
		Date & Time Collected <u>12/09/86</u>				Category _____	
AG	<10 ug/L	AS	28.7 ug/L	BE	<3 ug/L	CD	<5 ug/L
CU	11 ug/L	FE	20100 ug/L	HG	<0.2 ug/L	NI	<29 ug/L
SB	<44 ug/L	SE	<5 ug/L	TL	<10 ug/L	ZN	63 ug/L

CHECKLIST

SECTION II - Sample Data Packets

Sample No: SCK-P02-001

Organic Analysis Data Sheets (OADS)

VOA
 SV
 Pest.

Tentatively Identified Compounds

VOA
 SV

Raw Data

VOA chromatogram
 VOA quantitation report (surrogates)
 SV chromatogram
 SV quantitation report (surrogates)
 Pest. chromatogram and integrator print out

Inorganic Analysis Data Sheets (IADS)

003412

SPECTRIX DC # ----- 8

INST ID 400C

SAMPLE NUMBER: SCK-P02-001

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

LABORATORY NAME: SPECTRIX
 SAMPLE ID NO.: 861204107
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *[Signature]*

CASE NO.: -----
 QC REPORT NO.: -----
 CONTRACT NO.: -----
 DATE SAMPLE RECEIVED: 12/9/86

VOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED/PREPARED: 12/13/86
 DATE ANALYZED: 12/13/86

DATAFILE: 4U12041V07

DILUTION FACTOR: 20.0

CAS # COMPOUND

UG/L

003413

C010	CHLOROMETHANE	200	U
C015	BROMOMETHANE	200	U
C020	VINYL CHLORIDE	200	U
C025	CHLOROETHANE	200	U
C030	METHYLENE CHLORIDE	100	U
C035	ACETONE	200	U
C040	CARBON DISULFIDE	100	U
C045	1, 1-DICHLOROETHENE	100	U
C050	1, 1-DICHLOROETHANE	100	U
C055	TRANS-1, 2-DICHLOROETHENE	100	U
C060	CHLOROFORM	100	U
C065	1, 2-DICHLOROETHANE	100	U
C110	2-BUTANONE	200	U
C115	1, 1, 1-TRICHLOROETHANE	100	U
C120	CARBON TETRACHLORIDE	100	U
C125	VINYL ACETATE	200	U
C130	BROMODICHLOROMETHANE	100	U
C140	1, 2-DICHLOROPROPANE	100	U
C145	TRANS-1, 3-DICHLOROPROPENE	100	U
C150	TRICHLOROETHENE	100	U
C155	DIBROMOCHLOROMETHANE	100	U
C160	1, 1, 2-TRICHLOROETHANE	100	U
C165	BENZENE	70	J
C170	CIS-1, 3-DICHLOROPROPENE	100	U
C175	2-CHLOROETHYL VINYLETHER	200	U
C180	BROMOFORM	100	U
C210	2-HEXANONE	200	U
C205	4-METHYL-2-PENTANONE	200	U
C220	TETRACHLOROETHENE	100	U
C225	1, 1, 2, 2-TETRACHLOROETHANE	100	U
C230	TOLUENE	33	J
C235	CHLOROBENZENE	100	U
C240	ETHYLBENZENE	85	J
C245	STYRENE	100	U
C250	TOTAL XYLEMES	85	J

= UNDETECTED AT THE LISTED DETECTION LIMIT

= COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

INST ID: 4530

SPECTRIX DC # ---- 8
SAMPLE NUMBER: SCK-P02-001

ORGANICS ANALYSIS DATA SHEET - PAGE 2

LABORATORY NAME: SPECTRIX
 LAB SAMPLE ID NO.: 861204107
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *[Signature]*

CASE NO. : ---
 QC REPORT NO. : ---
 CONTRACT NO. : ---
 DATE SAMPLE RECEIVED: *12/9/86*

SEMIVOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED: *12/13/86*
 DATE ANALYZED: 01/11/87

DATAFILE: 4U12041C07A
 DILUTION FACTOR: *20*

CAS #	COMPOUND	UG/L	
C315	PHENOL	200	U
C325	BIS(2-CHLOROETHYL)ETHER	200	U
C330	2-CHLOROPHENOL	200	U
C335	1, 3-DICHLOROBENZENE	200	U
C340	1, 4-DICHLOROBENZENE	200	U
C345	BENZYL ALCOHOL	200	U
C350	1, 2-DICHLOROBENZENE	200	U
C355	2-METHYLPHENOL	200	U
C360	BIS(2-CHLOROISOPROPYL)ETHER	200	U
C365	4-METHYLPHENOL	200	U
C370	N-NITROSODIPIROPYLAMINE	200	U
C375	HEXAChLOROETHANE	200	U
C410	NITROBENZENE	200	U
C415	ISOPHORONE	200	U
C420	2-NITROPHENOL	200	U
C425	2, 4-DIMETHYLPHENOL	200	U
C430	BENZOIC ACID	1000	U
C435	BIS(2-CHLOROETHOXY)METHANE	200	U
C440	2, 4-DICHLOROPHENOL	200	U
C445	1, 2, 4-TRICHLOROBENZENE	200	U
C450	NAPHTHALENE	6100	
C455	4-CHLOROANILINE	200	U
C460	HEXAChLOROBUTADIENE	200	U
C465	P-CHLORO-M-CRESOL	200	U
C470	2-METHYLNAPHTHALENE	400	
C510	HEXAChLOROCYCLOPENTADIENE	200	U
C515	2, 4, 6-TRICHLOROPHENOL	200	U
C520	2, 4, 5-TRICHLOROPHENOL	1000	U
C525	2-CHLORONAPHTHALENE	200	U
C530	2-NITROANILINE	1000	U
C535	DIMETHYL PHTHALATE	200	U
C540	ACENAPHTHYLENE	17	J
C545	3-NITROANILINE	1000	U
C550	ACENAPHTHENE	1200	
C555	2, 4-DINITROPHENOL	1000	U
C560	4-NITROPHENOL	1000	U
C565	DIBENZOFURAN	620	
C570	2, 4-DINITROTOLUENE	200	U
C575	2, 6-DINITROTOLUENE	200	U

SPECTRIX DC # ---- 8

SAMPLE NUMBER: SCK-P02-001

MIVOLATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

DATAFILE: 4U12041C07A

CAS #	COMPOUND	UG/L
C580	DIETHYL PHTHALATE	200 U
C585	4-CHLOROPHENYL PHENYL ETHER	200 U
C590	FLUORENE	440
C595	4-NITROANILINE	1000 U
C610	4, 6-DINITRO-2-METHYLPHENOL	1000 U
C615	N-NITROSODIPHENYLAMINE	200 U
C625	4-BROMOPHENYL PHENYL ETHER	200 U
C630	HEXACHLOROBENZENE	200 U
C635	PENTACHLOROPHENOL	1000 U
C640	PHENANTHRENE	990
C645	ANTHRACENE	110 J
C650	DI-N-BUTYL PHTHALATE	200 U
C655	FLUORANTHENE	130 J
C715	PYRENE	140 J
C720	BUTYL BENZYL PHTHALATE	200 U
C725	3, 3'-DICHLOROBENZIDINE	400 U
C730	BENZO(A)ANTHRACENE	200 U
C735	BIS(2-ETHYLHEXYL)PHTHALATE	200 U
C740	CHRYSENE	200 U
C760	DI-N-OCTYL PHTHALATE	200 U
C765	BENZO(B)FLUORANTHENE	200 U
C770	BENZO(K)FLUORANTHENE	200 U
C775	BENZO(A)PYRENE	200 U
C780	INDENO(1, 2, 3-CD)PYRENE	200 U
C785	DIBENZO(A, H)ANTHRACENE	200 U
C790	BENZO(GHI)PERYLENE	200 U

003415

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

Lab Name: Spectrix
Case No: KEYSTONE
Lab No.: 86-12-041-07A

DC No: KEYSTONE
SMO No: SCK-P02-001

ORGANICS ANALYSIS DATA SHEET
(page 3)

Pesticide/PCBs

Conc: Low

GPC Cleanup Yes ✓
Sep. Funnel Ext. Yes
Cont. Liquid Ext. Yes

Date Extracted: 12/11/86

Date Analyzed: 12/24/86

Dilution Factor: 1 NA

Percent Moisture(decantated):

CAS #	Compound	ug/L
319-84-6	ALPHA-BHC	0.05 U
319-85-7	BETA-BHC	0.05 U
319-86-8	DELTA-BHC	0.05 U
58-89-9	GAMMA-BHC (LINDANE)	0.05 U
76-44-8	HEPTACHLOR	0.05 U
309-00-2	ALDRIN	0.05 U
1024-57-3	HEPTACHLOR EPOXIDE	0.05 U
959-98-8	ENDOSULFAN I	0.05 U
60-57-1	DIELDRIN	0.10 U
72-55-9	4,4'-DDE	0.10 U
72-20-8	ENDRIN	0.10 U
33213-65-9	ENDOSULFAN II	0.10 U
72-54-8	4,4'-DDD	0.10 U
7421-93-4	ENDRIN ALDEHYDE	0.10 U
1031-07-8	ENDOSULFAN SULFATE	0.10 U
50-29-3	4,4'-DDT	0.10 U
72-43-5	METHOXYSCHLOR	0.50 U
53494-70-5	ENDRIN KETONE	0.10 U
57-74-9	CHLORDANE	0.50 U
8001-35-2	TOXAPHENE	1.0 U
12674-11-2	AROCLOTR-1016	0.50 U
11104-28-2	AROCLOTR-1221	0.50 U
11141-16-5	AROCLOTR-1232	0.50 U
53469-21-9	AROCLOTR-1242	0.50 U
12672-29-6	AROCLOTR-1248	0.50 U
11097-69-1	AROCLOTR-1254	1.0 U
11096-82-5	AROCLOTR-1260	1.0 U

V_s = 500 = Volume of water extracted (ml)

V_t = 5000 = Volume of total extract (ul)

V_i = 1 = Volume of extract injected(ul)

1st Column Inst. : 6000

2nd Column Inst. : 6000 NA.

U=UNDETECTED AT THE LISTED DETECTION LIMIT.

J=COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT.

B=COMPOUND ALSO FOUND IN BLANK.

Form 1

INST ID: 4000

SPECTRIX DC # ----- 8

SAMPLE NUMBER: SCK-P02-001

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: -----

REPORT NO.: -----

ANALYST: CKT

DATAFILE: 4U12041V07

B. TENTATIVELY IDENTIFIED COMPOUNDS

AS #	VOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
124-38-9	CARBON DIOXIDE (ACN)*	43	945	420 ug/l
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	96		200 ug/l

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

* ALSO FOUND IN BLANK

INST ID: 4530

SPECTRIX DC # ---- 5

SAMPLE: SCK-P02-001

ORGANICS ANALYSIS DATA SHEET - PAGE 5

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: ---

GC REPORT NO.: ---

ANALYST: GU

DATAFILE: 4U12041C07A

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	SEMOVOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
700-41-4	BENZENE, ETHYL-			UG/L ⁰⁸
109-51-0	BENZENEACETIC ACID, 2-CARBOXY-	305	809	100 ³⁴
611-15-4	BENZENE, 1-ETHENYL-2-METHYL-	509	901	250 ^Q
125-13-6	1H-INDENE	562	958	480 ^J
95-15-8	BENZO[B]THIOPHENE	573	940	150 ^J
71-83-2	1H-INDENE, 1-ETHYLIDENE-	750	909	260 ^J
122-52-4	1, 1'-BIPHENYL	873	917	350 ^J
139-27-5	NAPHTHALENE, 2-ETHYL-	938	929	270 ^J
181-40-8	NAPHTHALENE, 2, 3-DIMETHYL-	952	901	87 ^J
581-40-8	NAPHTHALENE, 2, 3-DIMETHYL-	962	882	170 ^J
171-58-4	NAPHTHALENE, 1, 4-DIMETHYL-	976	935	300 ^J
126-53-3	1-NAPHTHALENECARBONITRILE	993	935	88 ^J
86-74-8	9H-CARBAZOLE	1042	787	120 ^J
		1313	894	290 ^J

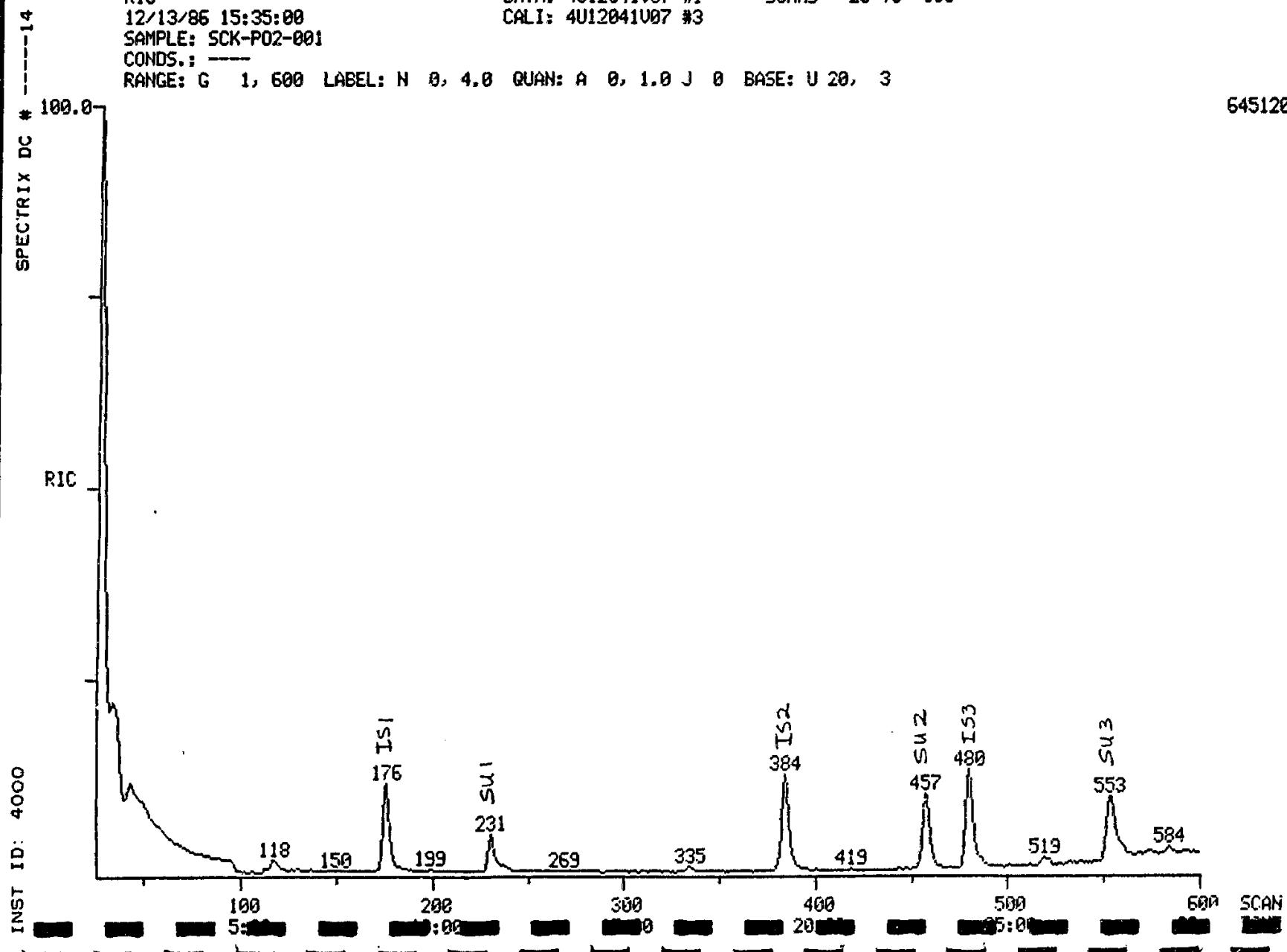
J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

003419

RIC
12/13/86 15:35:00
SAMPLE: SCK-P02-001
CONDNS.:
RANGE: G 1, 600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12041U07 #1 SCANS 25 TO 600
CALI: 4U12041U07 #3

645120.



SPECTRIX DC # -----14

VOA QUANTITATION REPORT

SAMPLE ID.	SCK-P02-001	INST	4000
FILENAME	4U12041V07	ANALYST	CKT
CLIENT	KEYSTONE	VERIFIED BY	<i>[Signature]</i>
DATE INJECTED	12/13/86 15:35:00	CORR. FACTOR	20.00
ALIB STD DATE	12-14-86		

CAN#	VOA COMPOUNDS	M/E	AMOUNT	RRT	AREA
175 C101	BROMOCHLOROMETHANE	**IS1**	128	1000 UG/L	1.000
384 C110	1, 4-DIFLUOROBENZENE	**IS2**	114	1000 UG/L	1.000
480 C120	D5-CHLOROBENZENE	**IS3**	117	1000 UG/L	1.000
231 CS15	D4-1, 2-DICHLOROETHANE	**SU1**	65	98 %	1.320
457 CS05	D8-TOLUENE	**SU2**	98	101 %	0.952
553 CS10	P-BROMOFLUOROBENZENE	**SU3**	95	93 %	1.152
117 C030	METHYLENE CHLORIDE *		84	129 UG/L	0.669
238 C110	2-BUTANONE *		72	221 UG/L	0.620
335 C165	BENZENE		78	70 UG/L	1.335
461 C230	TOLUENE		92	33 UG/L	0.872
519 C240	ETHYLBENZENE		106	85 UG/L	0.960
584 C250	TOTAL XYLEMES		106	85 UG/L	1.081
					2823
					3502

* BLANK CORRECTED ON DATA SHEET

003421

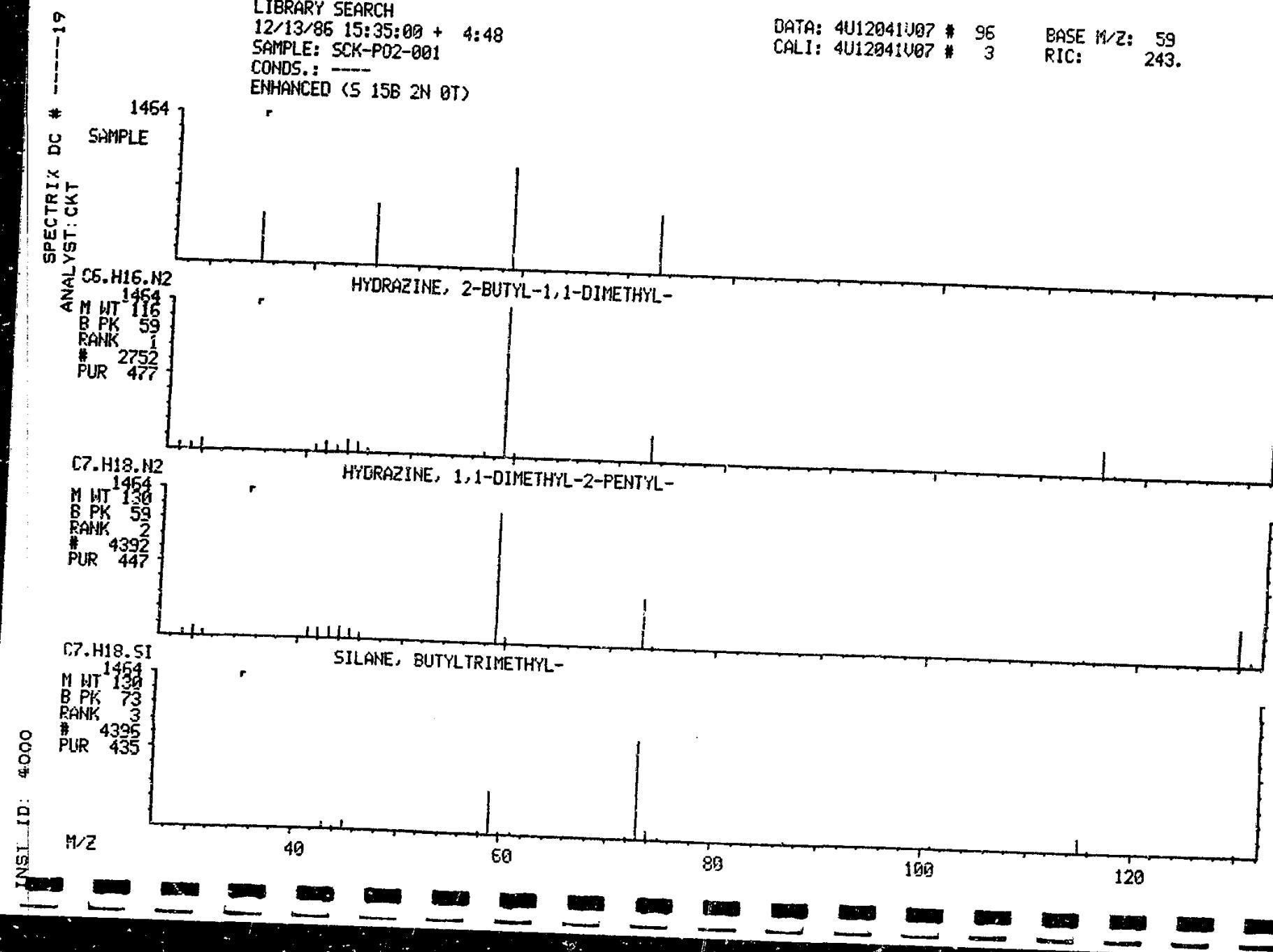
LIBRARY SEARCH

12/13/86 15:35:00 + 4:48

SAMPLE: SCK-P02-001

COND.: -----

ENHANCED (S 15B 2N 0T)

DATA: 4U12041U07 # 96
CALI: 4U12041U07 # 3BASE M/Z: 59
RIC: 243.

224200

RIC
01/11/87 13:55:00
SAMPLE: SCK-P02-001
COND.: ---
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 26, 3

DATA: 4U12041C07A #1 SCANS 200 TO 1350
CALI: 4U12041C07A #3

292624.

SPECTRIX DC # ----14

INST ID: 4530

100.0

RIC

200

31

305

406

463

624

527

562

793

816

873

938

976

1035

1118

1216

1274

1313

10:00

16:40

20

20:00

20

20:00

16:40

20:00

SCAN

TIME

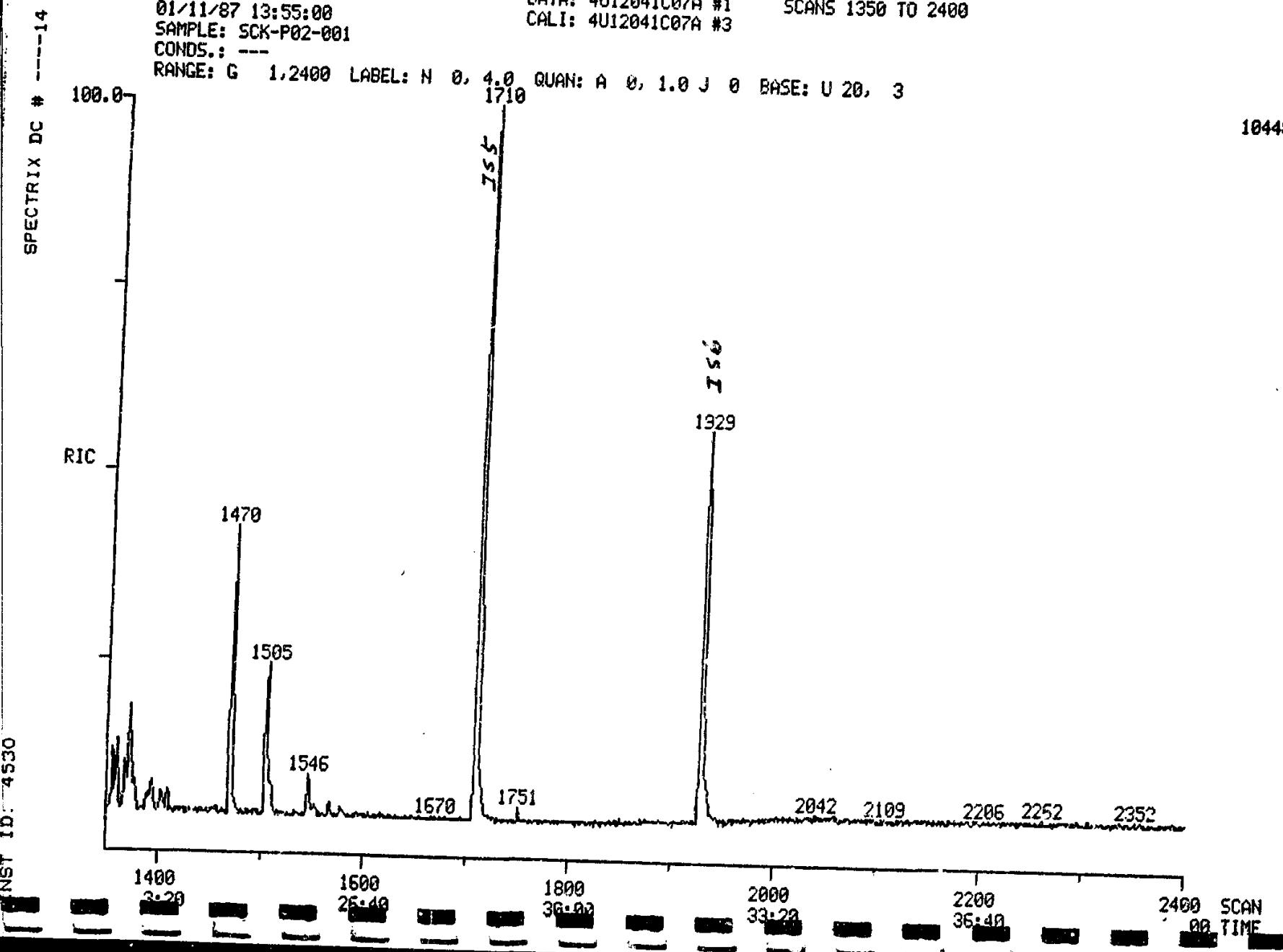
003422

003423

RIC
01/11/87 13:55:00
SAMPLE: SCK-P02-001
COND.: ---
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12041C07A #1
CALI: 4U12041C07A #3
SCANS 1350 TO 2400

10448.



003423

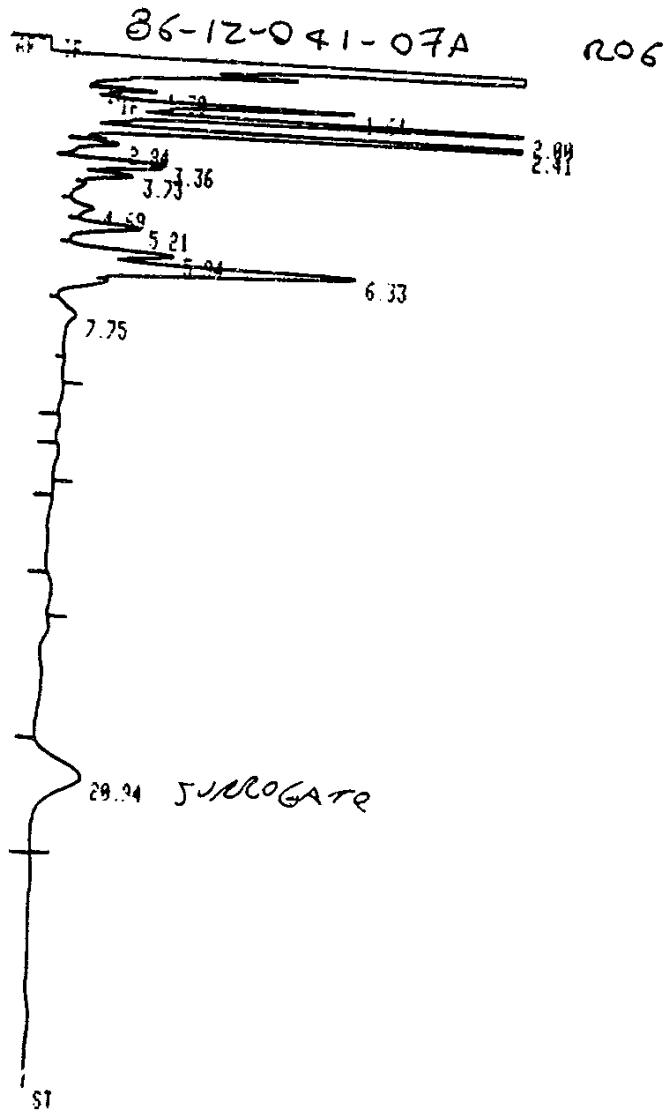
SPECTRIX DC # ---14

SEMIVOLATILE WATER REPORT

SAMPLE ID.	SCK-P02-001		
FILENAME	4U12041C07A	INST ID.	4530
CLIENT	KEYSTONE	ANALYST	GU
DATE INJECTED	01/11/87 13:55:00	VERIFIED BY	<i>YPM</i>
CALIB STD DATE	01/08/87	CORR. FACTOR	20.00

SCAN#	SEMIVOLATILE COMPOUNDS		M/E	AMOUNT	RRT	AREA
527	C130	1, 4-DICHLOROBENZEND-D4 ** IS1 **	152	800 UG/L	1.000	20829
738	C140	NAPHTHALENE-D8 ** IS2 **	136	800 UG/L	1.000	78449
1029	C150	ACENAPHTHENE-D10 ** IS3 **	164	800 UG/L	1.000	36511
1270	C160	PHENANTHRENE-D10 ** IS4 **	188	800 UG/L	1.000	40427
1710	C170	CHRYSENE-D12 ** IS5 **	240	800 UG/L	1.000	16890
1929	C175	PERYLENE-D12 ** IS6 **	264	800 UG/L	1.000	10686
744	C430	NAPHTHALENE	128	6150 UG/L	1.008	577467
856	C470	2-METHYLNAPHTHALENE	142	405 UG/L	1.160	35600
1002	C540	ACENAPHTHYLENE	152	17 UG/L	0.974	1230
1035	C550	ACENAPHTHENE	153	1191 UG/L	1.006	50021
1061	C565	DIBENZOFURAN	168	622 UG/L	1.031	35680
1118	C590	FLUORENE	166	443 UG/L	1.086	21068
1274	C640	PHENANTHRENE	178	987 UG/L	1.003	48019
1280	C645	ANTHRACENE	178	109 UG/L	1.008	5275
470	C655	FLUORANTHENE	202	132 UG/L	1.157	6065
505	C715	PYRENE	202	142 UG/L	0.880	3779

* SURROGATES WERE DILUTED OUT



DC # N/A

SPECTRIX CORPORATION

CASE NO. Keystone

SAMPLE NO. SEK-PZ-001

PESTICIDE FRACTION

1st Column - ID

1.8m X 4mm ID X 6mm 00

1.5% SP 2250

1.95% SP2401

100/120 Supelcoport

Column Temp. 203 °C

Volume Inj. 1.0 ul

Instrument ID VARIAN 600

RUN # 143
WORKFILE ID: C
WORKFILE NAME:

DEC/24/86 00:26:44

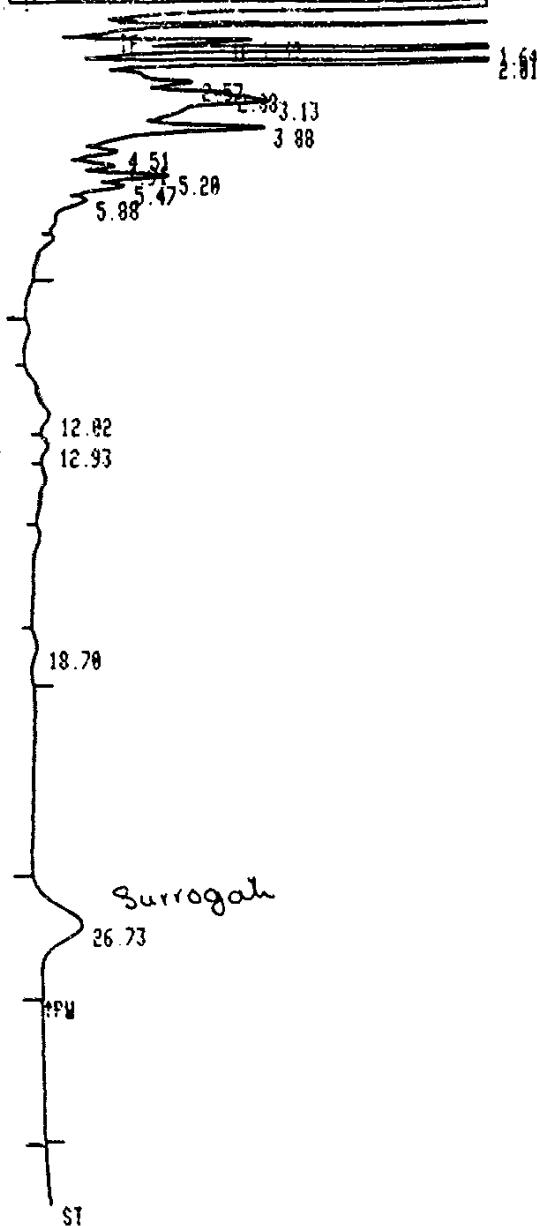
HEIGHT%

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
1.30	7695	D BY	0.113	2.524
1.64	33853	VV	0.166	11.236
2.00	67261	D VV	0.122	22.324
2.41	92318	VB	0.121	30.640
2.84	4827	BP	0.165	1.602
3.36	12460	PY	0.136	4.135
3.73	8553	VV	0.182	2.839
4.69	3765	VV	0.367	1.250
5.21	9745	VV	0.289	3.234
5.94	14509	VV	0.259	4.915
6.33	39208	VV	0.290	12.681
7.25	2184	PY	0.622	0.725
20.94	6013	PB	1.006	1.996

TOTAL HGT= 301300
MUL FACTOR= 1.0000E+00

START: 86-12-041-07A

NA.



DC# _____
SPECTRIX HOUSTON

CASE NO. Keystone

SAMPLE NO. Sck-poz-0c
PESTICIDE FRACTION
2ND COLUMN CONFIRMATI

2ND COLUMN TLC
1.8m x2mm ID x6mm OD
3% SP-2100
100/120 SUPERLCOPORT
COLUMN TEMP. 187 °C
VOLUME INJ. 1 μl N
INSTRUMENT ID. Gooch

M
O
O

RUN # 15 JAN/06/87 19:38:14

WORKFILE ID: C

WORKFILE NAME:

HEIGHT%	RT	HEIGHT	TYPE	AR/HT	HEIGHT%
	1.39	43712	PY	0.166	4.963
	1.64	323694	D VV	0.099	36.752
	2.01	230495	D VV	0.107	26.170
	2.57	30379	VV	0.348	3.449
	2.88	40327	VV	0.173	4.579
	3.13	51507	VV	0.508	5.848
	3.88	51704	VV	0.334	5.871
	4.51	14656	VV	0.260	1.664
	4.91	14780	VV	0.230	1.678
	5.20	28862	VV	0.235	3.277
	5.47	18000	VV	0.277	2.044
	5.88	9138	VV	0.475	1.038
	12.02	5593	PY	1.124	0.635
	12.93	5134	VV	0.681	0.583
	18.70	1300	VB	0.542	0.148
	26.73	11463	BB	1.182	1.302

Page 1

Received: 12/09/86

SPECTRUM 200 REPORT
Results by Sample

Work Order # 86-12-041

SAMPLE ID <u>SCK-PD2-001 (LLW)</u>			SAMPLE # <u>Q7</u> FRACTIONS: <u>A</u>								
			Date & Time Collected <u>12/09/86</u>			Category _____					
AG	<10	AS	29.5	BE	<3	CD	<5	CN_SU	<100	CR	<9
CU	<10	FE	7380	HG	<0.2	NI	64	N03_SU	<100	PB	<5
SB	<44	SE	<5	TL	<10	ZN	33				

CHECKLIST

SECTION II - Sample Data Packets

Sample No: Trip Blank

Organic Analysis Data Sheets (OADS)

VOA
 N/A SV
 N/A Pest.

Tentatively Identified Compounds

VOA
 N/A SV

Raw Data

VOA chromatogram
 VOA quantitation report (surrogates)
 N/A SV chromatogram
 N/A SV quantitation report (surrogates)
 N/A Pest. chromatogram and integrator print out

N/A Inorganic Analysis Data Sheets (IADS)

003428

SPECTRIX DC # ----- 8

INST ID 4000

SAMPLE NUMBER: TRIP BLANK

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

LABORATORY NAME: SPECTRIX

CASE NO.: ----

SAMPLE ID NO.: 861204106

GC REPORT NO.: ----

SAMPLE MATRIX: WATER

CONTRACT NO.: ----

DATA RELEASE AUTHORIZED BY: *MTH*

DATE SAMPLE RECEIVED: 12/4/86

VOLATILES

CONCENTRATION: LOW

DATAFILE: 4U12041V06

DATE EXTRACTED/PREPARED: 12/13/86

DILUTION FACTOR: 10.....

DATE ANALYZED: 12/13/86

CAS #	COMPOUND	UG/L	DET
C010	CHLOROMETHANE	10 U	00342
C015	BROMOMETHANE	10 U	
C020	VINYL CHLORIDE	10 U	
C025	CHLOROETHANE	10 U	
C030	METHYLENE CHLORIDE	5 U	
C035	ACETONE	10 U	
C040	CARBON DISULFIDE	5 U	
C045	1, 1-DICHLOROETHENE	5 U	
C050	1, 1-DICHLOROETHANE	5 U	
C055	TRANS-1, 2-DICHLOROETHENE	5 U	
C060	CHLOROFORM	5 U	
C065	1, 2-DICHLOROETHANE	5 U	
C110	2-BUTANONE	10 U	
C115	1, 1, 1-TRICHLOROETHANE	5 U	
C120	CARBON TETRACHLORIDE	5 U	
C125	VINYL ACETATE	10 U	
C130	BROMODICHLOROMETHANE	5 U	
C140	1, 2-DICHLOROPROPANE	5 U	
C145	TRANS-1, 3-DICHLOROPROPENE	5 U	
C150	TRICHLOROETHENE	5 U	
C155	DIBROMOCHLOROMETHANE	5 U	
C160	1, 1, 2-TRICHLOROETHANE	5 U	
C165	BENZENE	5 U	
C170	CIS-1, 3-DICHLOROPROPENE	5 U	
C175	2-CHLOROETHYL VINYLETHER	10 U	
C180	BROMOFORM	5 U	
C210	2-HEXANONE	10 U	
C205	4-METHYL-2-PENTANONE	10 U	
C220	TETRACHLOROETHENE	5 U	
C225	1, 1, 2, 2-TETRACHLOROETHANE	5 U	
C230	TOLUENE	5 U	
C235	CHLOROBENZENE	5 U	
C240	ETHYLBENZENE	5 U	
C245	STYRENE	5 U	
C250	TOTAL XYLENES	5 U	

= UNDETECTED AT THE LISTED DETECTION LIMIT

= COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

INST-ID: 4000

SPECTRIX DC # ----- 8

SAMPLE NUMBER: TRIP BLANK

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: -----

LC REPORT NO.: -----

ANALYST: CKT

DATAFILE: 4U12041V06

B. TENTATIVELY IDENTIFIED COMPOUNDS

AS #	VOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
124-38-9	CARBON DIOXIDE (ACN) *	34	967	15 4 3
124-38-9	CARBON DIOXIDE (ACN) *	49	966	18 0 3
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	96		7 0 0

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

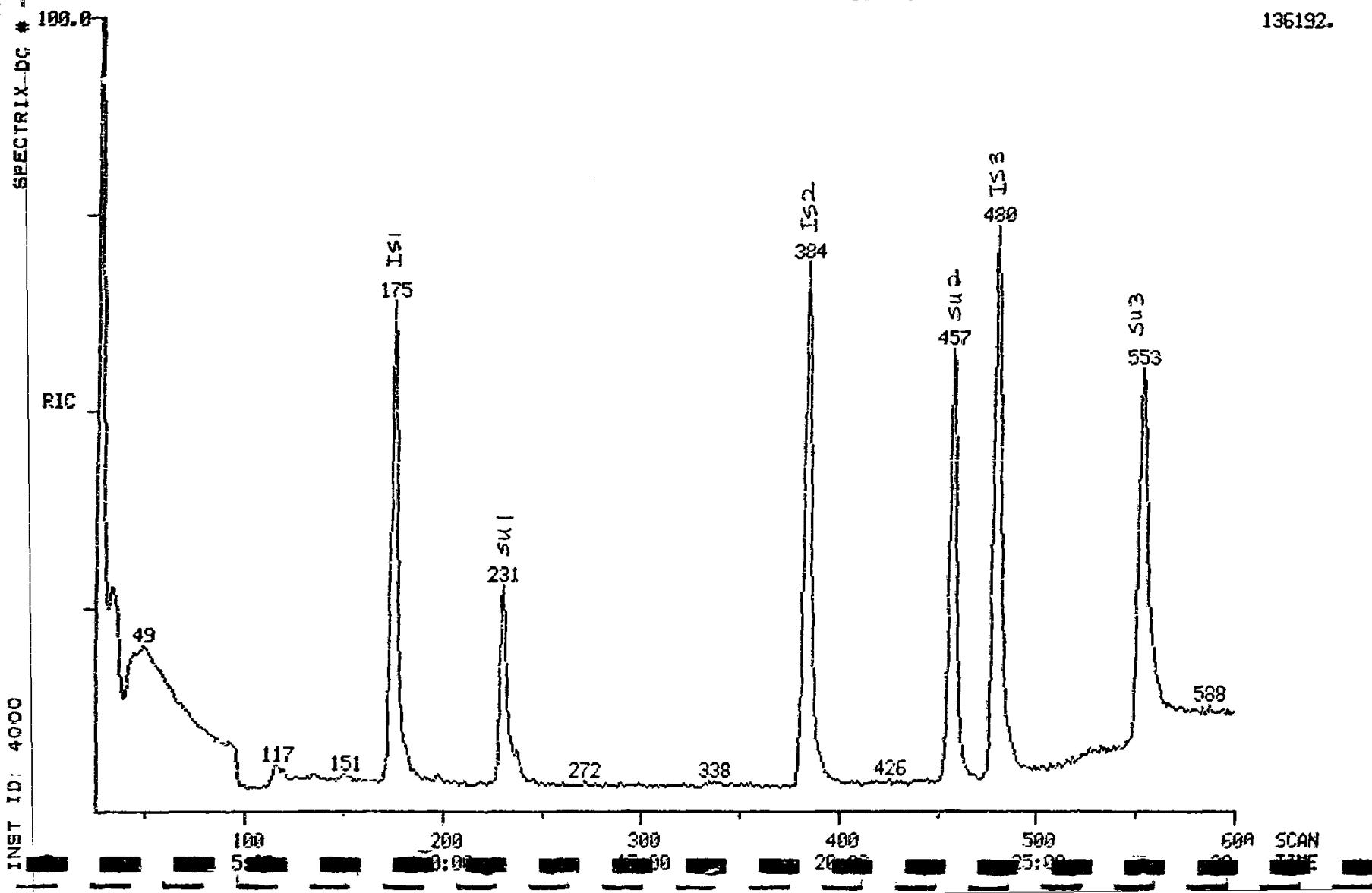
* ALSO FOUND IN BLANK

003431

RIC
12/13/86 14:10:00
SAMPLE: TRIP BLANK
COND.:
RANGE: G 1, 600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4U12041V06 #1 SCANS 25 TO 600
CALI: 4U12041V06 #3

136192.



003431

SPECTRIX DC #-----14

VOA QUANTITATION REPORT

SAMPLE ID.	TRIP BLANK		
JENAME	4U12041V06	INST	4000
CLIENT	KEYSTONE	ANALYST	CKT
DATE INJECTED	12/13/86 14:10:00	VERIFIED BY	<i>SPH</i>
ALIB STD DATE	12-14-86	CORR. FACTOR	1.00

CAN#	VOA COMPOUNDS	M/E	AMOUNT	RRT	AREA
175 C101	BROMOCHLOROMETHANE	**IS1**	128	50 UG/L	1.000 35698
384 C110	1, 4-DIFLUOROBENZENE	**IS2**	114	50 UG/L	1.000 144699
480 C120	D5-CHLOROBENZENE	**IS3**	117	50 UG/L	1.000 116200
231 CS15	D4-1, 2-DICHLOROETHANE	**SU1**	65	96 %	1.320 35225
457 CS05	D8-TOLUENE	**SU2**	98	104 %	0.952 99739
553 CS10	P-BROMOFLUOROBENZENE	**SU3**	95	100 %	1.152 60454
237 C110	2-BUTANONE *		72	12 UG/L	0.617 1745

* BLANK CORRECTED ON DATA SHEET

00343

003433

19

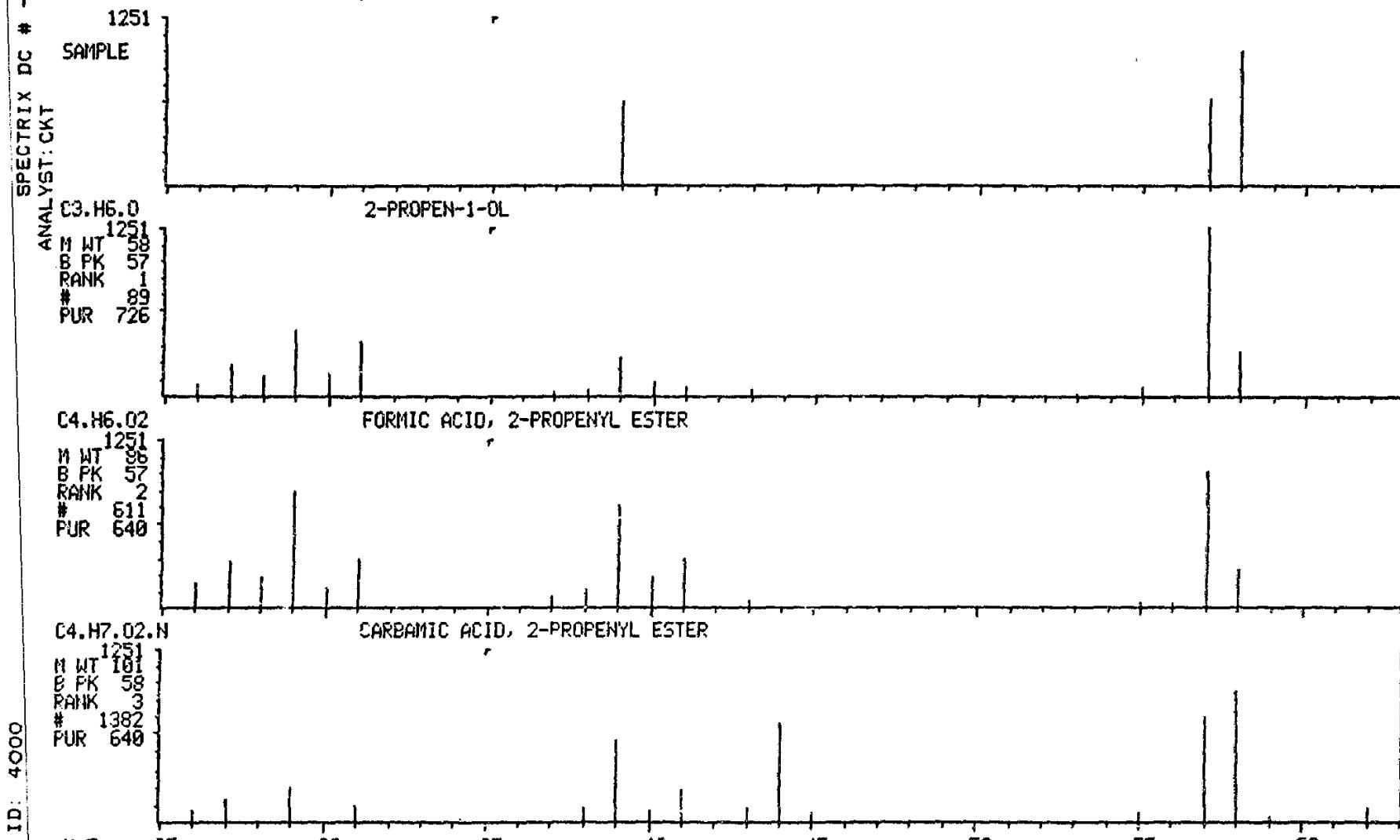
SPECTRUM DC #

ANALYST: CKT

INST ID: 4000

LIBRARY SEARCH
12/13/86 14:10:00 + 4:48
SAMPLE: TRIP BLANK
COND'S: ----
ENHANCED (S 15B 2N 0T)

DATA: 4U12041V06 # 96
CALI: 4U12041V06 # 3
BASE M/Z: 58
RIC: 132.



CHECKLIST

SECTION III - Blanks

OADS

VOA
 SV
 Pest.

Tentatively Identified Compounds

VOA
 SV

Raw Data

VOA chromatogram
 VOA quantitation report (surrogates)
 SV chromatogram
 SV quantitation report (surrogates)
 Pest. chromatogram with integrator print out

IADS

003434

SPECTRIX DC# ---- 8

NET ID 4000

SAMPLE NUMBER: LAB BLANK

ORGANICS ANALYSIS DATA SHEET (PAGE 1)

LABORATORY NAME: SPECTRIX

CASE NO.: ----

SAMPLE ID NO.: 8612041

GC REPORT NO.: ----

SAMPLE MATRIX: WATER

CONTRACT NO.: ----

DATA RELEASE AUTHORIZED BY: *ypm*

DATE SAMPLE RECEIVED: 12/9/86

VOLATILES

CONCENTRATION: LOW

DATAFILE: 4B1213V01

DATE EXTRACTED/PREPARED: 12/13/86

DILUTION FACTOR: 10.....

DATE ANALYZED: 12/13/86

CAS # COMPOUND

UG/L

003435

C010	CHLOROMETHANE	10	U
C015	BROMOMETHANE	10	U
C020	VINYL CHLORIDE	10	U
C025	CHLOROETHANE	10	U
C030	METHYLENE CHLORIDE	5	U
C035	ACETONE	10	U
C040	CARBON DISULFIDE	5	U
C045	1, 1-DICHLOROETHENE	5	U
C050	1, 1-DICHLOROETHANE	5	U
C055	TRANS-1, 2-DICHLOROETHENE	5	U
C060	CHLOROFORM	5	U
C065	1, 2-DICHLOROETHANE	5	U
C110	2-BUTANONE	10	U
C115	1, 1, 1-TRICHLOROETHANE	5	U
C120	CARBON TETRACHLORIDE	5	U
C125	VINYL ACETATE	10	U
C130	BROMODICHLOROMETHANE	5	U
C140	1, 2-DICHLOROPROPANE	5	U
C145	TRANS-1, 3-DICHLOROPROPENE	5	U
C150	TRICHLOROETHENE	5	U
C155	DIBROMOCHLOROMETHANE	5	U
C160	1, 1, 2-TRICHLOROETHANE	5	U
C165	BENZENE	5	U
C170	CIS-1, 3-DICHLOROPROPENE	5	U
C175	Z-CHLOROETHYL VINYLETHER	10	U
C180	BROMOFORM	5	U
C210	2-HEXANONE	10	U
C205	4-METHYL-2-PENTANONE	10	U
C220	TETRACHLOROETHENE	5	U
C225	1, 1, 2, 2-TETRACHLOROETHANE	5	U
C230	TOLUENE	5	U
C235	CHLOROBENZENE	5	U
C240	ETHYL BENZENE	5	U
C245	STYRENE	5	U
C250	TOTAL XYLENES	5	U

= UNDETECTED AT THE LISTED DETECTION LIMIT

= COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

= COMPOUND ALSO FOUND IN BLANK

INST ID 4530

SPECTRIX DC # ---- 8
SAMPLE NUMBER: LLW BLK

ORGANICS ANALYSIS DATA SHEET - PAGE 2

LABORATORY NAME: SPECTRIX
 AB SAMPLE ID NO.: B612041
 SAMPLE MATRIX: WATER
 DATA RELEASE AUTHORIZED BY: *[Signature]*

CASE NO.: ---
 QC REPORT NO.: ---
 CONTRACT NO.: ---
 DATE SAMPLE RECEIVED: 12/19/86

SEMIVOLATILES

CONCENTRATION: LOW
 DATE EXTRACTED: 12/13/86
 DATE ANALYZED: 01/09/87

DATAFILE: 4B12041C01

DILUTION FACTOR: 1.0

CAS #	COMPOUND	UG/L	
C315	PHENOL	10 U	6
C325	BIS(2-CHLOROETHYL)ETHER	10 U	M
C330	2-CHLOROPHENOL	10 U	4
C335	1, 3-DICHLOROBENZENE	10 U	N
C340	1, 4-DICHLOROBENZENE	10 U	0
C345	BENZYL ALCOHOL	10 U	0
C350	1, 2-DICHLOROBENZENE	10 U	0
C355	2-METHYLPHENOL	10 U	0
C360	BIS(2-CHLOROISOPROPYL)ETHER	10 U	0
C365	4-METHYLPHENOL	10 U	0
C370	N-NITROSODIPROPYLAMINE	10 U	0
C375	HEXAChLORoETHANE	10 U	0
C410	NITROBENZENE	10 U	0
C415	ISOPHORONE	10 U	0
C420	2-NITROPHENOL	10 U	0
C425	2, 4-DIMETHYLPHENOL	10 U	0
C430	BENZOIC ACID	50 U	0
C435	BIS(2-CHLOROETHOXY)METHANE	10 U	0
C440	2, 4-DICHLOROPHENOL	10 U	0
C445	1, 2, 4-TRICHLOROBENZENE	10 U	0
C450	NAPHTHALENE	10 J	0
C455	4-CHLORANILINE	10 U	0
C460	HEXAChLOROBUTADIENE	10 U	0
C465	P-CHLORO-M-CRESOL	10 U	0
C470	2-METHYLNAPHTHALENE	10 U	0
C510	HEXAChLOROCYCLOPENTADIENE	10 U	0
C515	2, 4, 6-TRICHLOROPHENOL	10 U	0
C520	2, 4, 5-TRICHLOROPHENOL	50 U	0
C525	2-CHLORONAPHTHALENE	10 U	0
C530	2-NITROANILINE	50 U	0
C535	DIMETHYL PHTHALATE	10 U	0
C540	ACENAPHTHYLENE	10 U	0
C545	3-NITROANILINE	50 U	0
C550	ACENAPHTHENE	10 U	0
C555	2, 4-DINITROPHENOL	50 U	0
C560	4-NITROPHENOL	50 U	0
C565	DIBENZOFURAN	10 U	0
C570	2, 4-DINITROTOLUENE	10 U	0
C575	2, 6-DINITROTOLUENE	10 U	0

SPECTRIX DC # ---- 8

SAMPLE NUMBER: LLW BLK

EMIVOLATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

DATAFILE: 4B12041C01

CAS #	COMPOUND	UG/L
C580	DIETHYL PHTHALATE	1 J
C585	4-CHLOROPHENYL PHENYL ETHER	10 U
C590	FLUORENE	10 U
C595	4-NITROANILINE	50 U
C610	4, 6-DINITRO-2-METHYLPHENOL	50 U
C615	N-NITROSODIPHENYLAMINE	10 U
C625	4-BROMOPHENYL PHENYL ETHER	10 U
C630	HEXACHLOROBENZENE	10 U
C635	PENTACHLOROPHENOL	50 U
C640	PHENANTHRENE	10 U
C645	ANTHRACENE	10 U
C650	DI-N-BUTYL PHTHALATE	10 U
C655	FLUORANTHENE	10 U
C715	PYRENE	10 U
C720	BUTYL BENZYL PHTHALATE	10 U
C725	3, 3'-DICHLOROBENZIDINE	20 U
C730	BENZO(A)ANTHRACENE	10 U
C735	BIS(2-ETHYLHEXYL)PHTHALATE	10 U
C740	CHRYSENE	10 U
C760	DI-N-OCTYL PHTHALATE	10 U
C765	BENZO(B)FLUORANTHENE	10 U
C770	BENZO(K)FLUORANTHENE	10 U
C775	BENZO(A)PYRENE	10 U
C780	INDENO(1, 2, 3-CD)PYRENE	10 U
C785	DIBENZO(A, H)ANTHRACENE	10 U
C790	BENZO(GHI)PERYLENE	10 U

J = UNDETECTED AT THE LISTED DETECTION LIMIT

U = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

003437

Lab Name: Spectrix
Case No: KEYSTONE
Lab No.: 86-12-041-01A

DC No: KEYSTONE
SMD No: L.BLANK LLW

ORGANICS ANALYSIS DATA SHEET
(page 3)

Pesticide/PCBs

Conc: Low
Date Extracted: 12/11/86
Date Analyzed: 12/23/86
Dilution Factor: 1
Percent Moisture(decanted):

GPC Cleanup Yes ✓ No
Sep. Funnel Ext. Yes
Cont. Liquid Ext. Yes ✓

CAS #	Compound	ug/L
319-84-6	ALPHA-BHC	0.05 U
319-85-7	BETA-BHC	0.05 U
319-86-8	DELTA-BHC	0.05 U
58-89-9	GAMMA-BHC (LINDANE)	0.05 U
76-44-8	HEPTACHLOR	0.05 U
309-00-2	ALDRIN	0.05 U
1024-57-3	HEPTACHLOR EPOXIDE	0.05 U
959-98-8	ENDOSULFAN I	0.05 U
60-57-1	DIELDRIN	0.10 U
72-55-9	4,4'-DDE	0.10 U
72-20-8	ENDRIN	0.10 U
33213-65-9	ENDOSULFAN II	0.10 U
72-54-8	4,4'-DDD	0.10 U
7421-93-4	ENDRIN ALDEHYDE	0.10 U
1031-07-8	ENDOSULFAN SULFATE	0.10 U
50-29-3	4,4'-DDT	0.10 U
72-43-5	METHOXYCHLOR	0.50 U
53494-70-5	ENDRIN KETONE	0.10 U
57-74-9	CHLORDANE	0.50 U
8001-35-2	TUXAPHENE	1.0 U
12674-11-2	AROCLOR-1016	0.50 U
11104-28-2	AROCLOR-1221	0.50 U
11141-16-5	AROCLOR-1232	0.50 U
53469-21-9	AROCLOR-1242	0.50 U
12672-29-6	AROCLOR-1248	0.50 U
11097-69-1	AROCLOR-1254	1.0 U
11096-82-5	AROCLOR-1260	1.0 U

V_s = 500 = Volume of water extracted (ml)

V_t = 5000 = Volume of total extract (ul)

V_i = 1 = Volume of extract injected(ul)

1st Column Inst. : 6000

2nd Column Inst. : 6000 NA.

U=UNDETECTED AT THE LISTED DETECTION LIMIT.

J=COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT.

B=COMPOUND ALSO FOUND IN BLANK.

INST ID: 4000

SPECTRIX DC # ----- 8

SAMPLE NUMBER: LAB BLANK

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: -----

QC REPORT NO.: -----

ANALYST: CKT

DATAFILE: 4B1213V01

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	VOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
24-38-9	CARBON DIOXIDE (ACN)	48	94%	100%
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	96	9%	10%

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

INST ID: 4530

SPECTRIX DC # ---- 8

SAMPLE: LLW BLK

ORGANICS ANALYSIS DATA SHEET - PAGE 5

LABORATORY NAME: SPECTRIX CORPORATION

CASE NO.: ---

GC REPORT NO.: ---

ANALYST: GU

DATAFILE: 4B12041C01

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	SEMOVOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
110-12-3	2-HEXANONE, 5-METHYL-	333	927	62 4
140-97-6	CYCLOHEXASILOXANE, DODECAMETHYL-	882	854	5 0
540-97-6	CYCLOHEXASILOXANE, DODECAMETHYL-	895	843	4 0
	UNKNOWN-DOESN'T MATCH ANY LIBRARY SPECTRA	1047		7 J
71890-38-2	9-EICOSYNE	1464	805	73 J
2075-51-4	1, 12-DODECANEDIOL	1502	770	6 J

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

003441

RIC
12/13/86 10:40:00
SAMPLE: LAB BLANK
COND.: ---
RANGE: G 1, 600 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4B1213U01 #1 SCANS 30 TO 600
CALI: 4B1213U01 #3

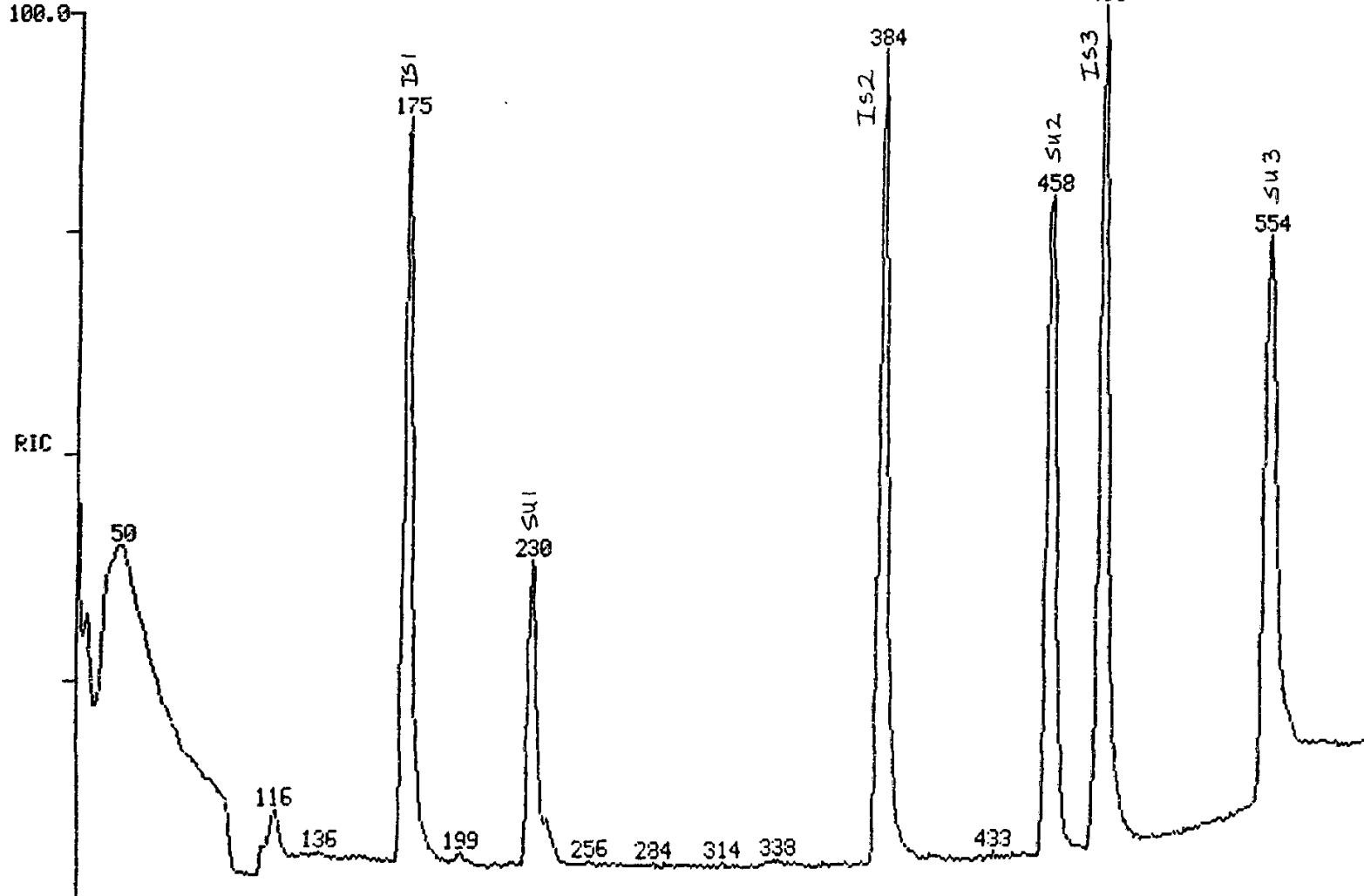
SPECTRIX DC # -----14

100.0

RIC

INST ID: 4000

003441



83328.

SPECTRIX DC # -----14

VOA QUANTITATION REPORT

SAMPLE ID. LAB BLANK
 FILENAME 4B1213V01
 CLIENT KEYSTONE INST 4000
 DATE INJECTED 12/13/86 10:40:00 ANALYST CKT
 CALIB STD DATE 12-14-86 VERIFIED BY *YTH*
 CORR. FACTOR 1.00

SCAN#	VOA COMPOUNDS		M/E	AMOUNT	RRT	AREA
174 CI01	BROMOCHLOROMETHANE	**IS1**	128	50 UG/L	1.000	30379
384 CI10	1, 4-DIFLUOROBENZENE	**IS2**	114	50 UG/L	1.000	125463
180 CI20	D5-CHLOROBENZENE	**IS3**	117	50 UG/L	1.000	100443
230 CS15	D4-1, 2-DICHLOROETHANE	**SU1**	65	96 %	1.322	298050
457 CS05	D8-TOLUENE	**SU2**	98	99 %	0.952	82044
553 CS10	P-BROMOFLUOROBENZENE	**SU3**	95	99 %	1.152	49680
116 C030	METHYLENE CHLORIDE		84	3 UG/L	0.667	2385
237 C110	2-BUTANONE		72	10 UG/L	0.617	1211

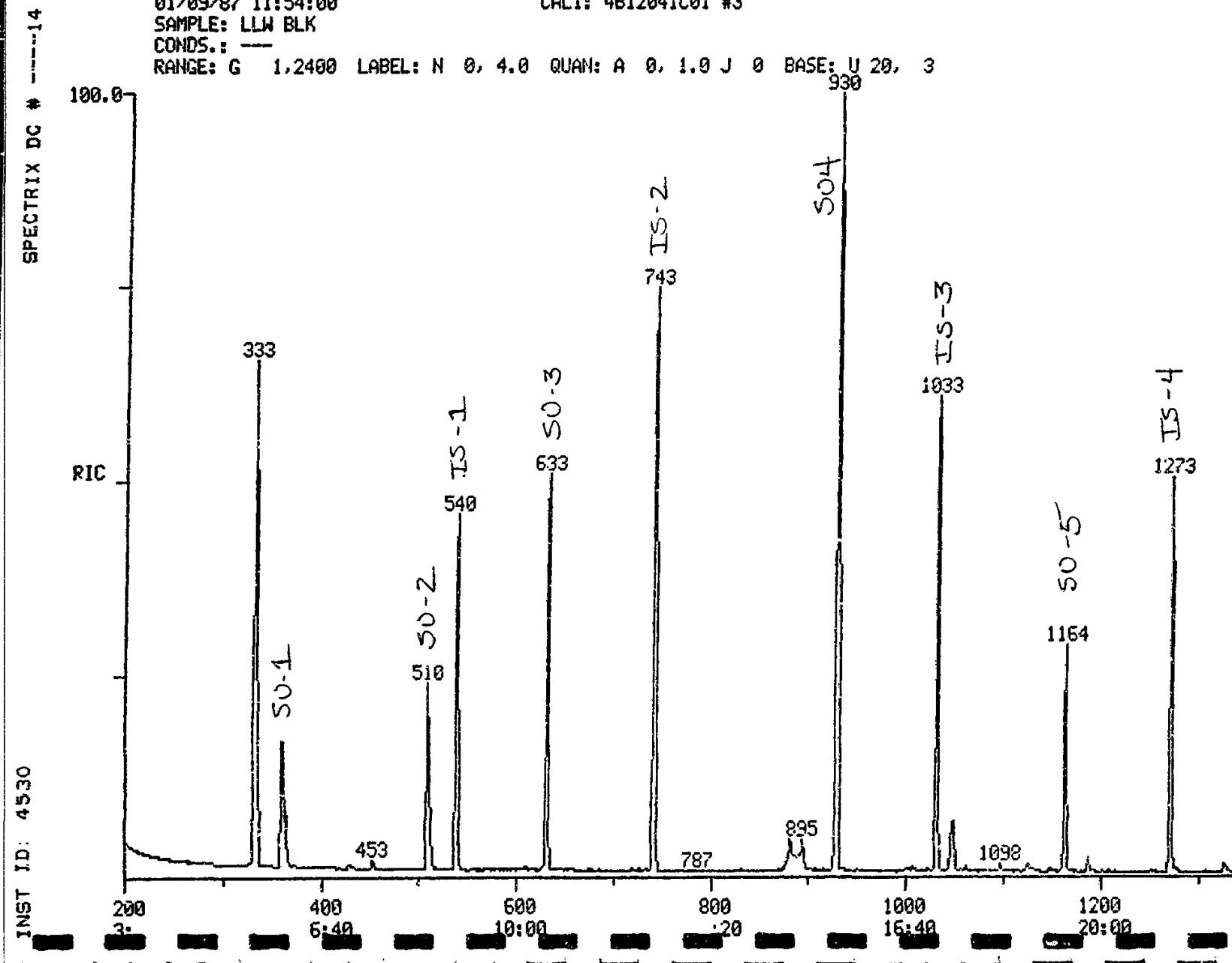
00344

003443

RIC
01/09/87 11:54:00
SAMPLE: LLW BLK
COND.: ---
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4B12041C01 #1 SCANS 200 TO 1350
CALI: 4B12041C01 #3

SPECTRIX DC # -----14



175104.

003443

003444

RIC
01/09/87 11:54:00
SAMPLE: LLW BLK
COND.: —
RANGE: G 1,2400 LABEL: N 0, 4.0 QUAN: A 0, 1.0 J 0 BASE: U 20, 3

DATA: 4B12041C01 #1 SCANS 1350 TO 2400
CALI: 4B12041C01 #3

144128.

SPECTRIX DC # -----14

RIC

INST ID: 4550

100.0

1400
3:20

1600
26:40

1800
30:00

2000
33:20

2200
36:40

2400
1:00 SCAN
TIME

1397

1590 1645

1753 1830

1975

2107 2159

2254 2347

SO-6

1549

TS-5

1712

TS-6

1932

003444

SPECTRIX DC # ----14

SEMIVOLATILE WATER REPORT

SAMPLE ID.

LLW BLK

FILENAME

4B12041C01

INST ID.

4530

CLIENT

KEYSTONE

ANALYST

GU

DATE INJECTED

01/09/87 11:54:00

VERIFIED BY

CALIB STD DATE

01/08/87

CORR. FACTOR

1.00

SCAN#

SEMIVOLATILE COMPOUNDS

M/E

AMOUNT

RRT

AREA

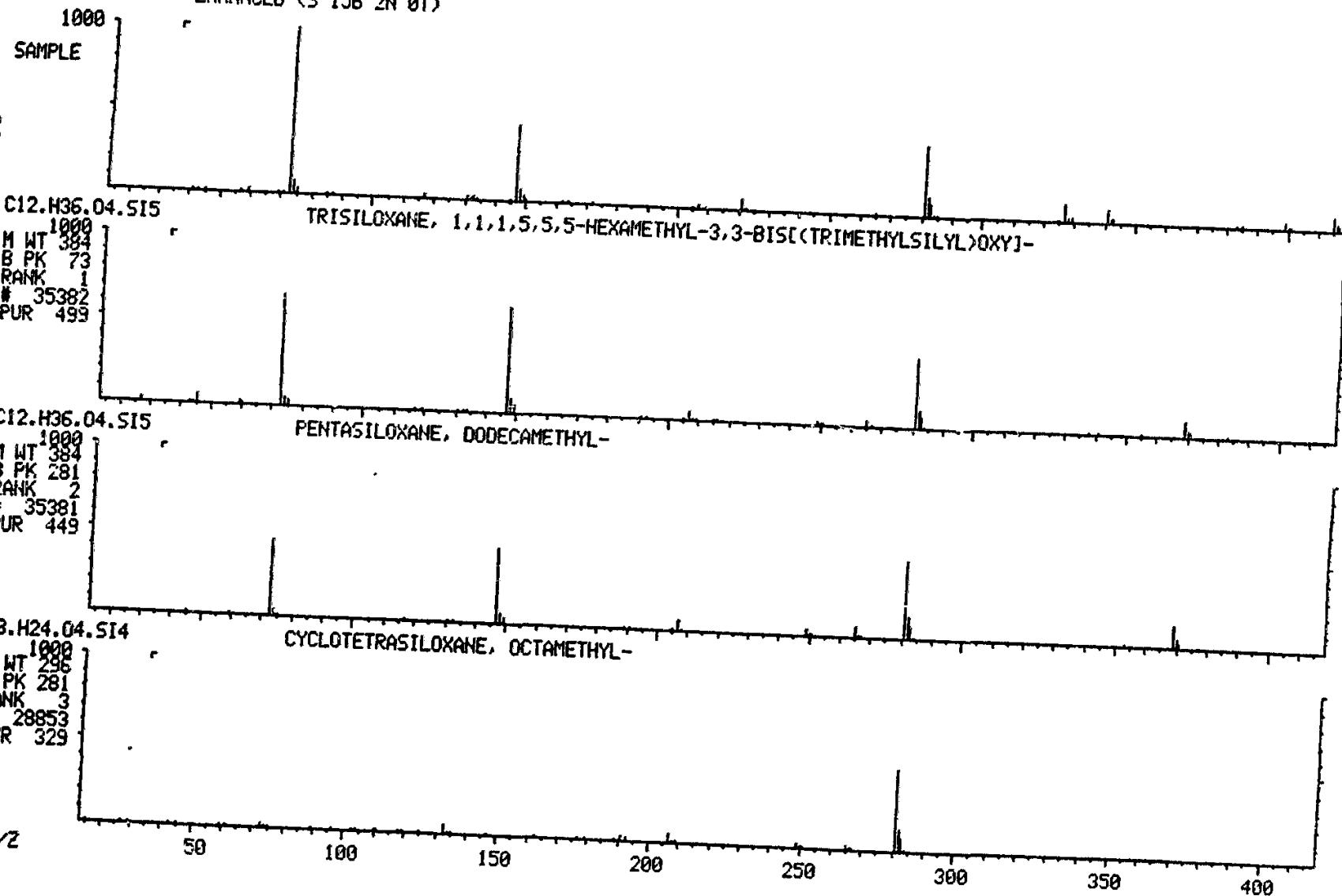
540	CI30	1,4-DICHLOROBENZEND-D4 ** IS1 **	152	40 UG/L	1.000	3599:
743	CI40	NAPHTHALENE-D8 ** IS2 **	136	40 UG/L	1.000	14708:
1033	CI50	ACENAPHTHENE-D10 ** IS3 **	164	40 UG/L	1.000	7625:
1273	CI60	PHENANTHRENE-D10 ** IS4 **	188	40 UG/L	1.000	9905:
1712	CI70	CHRYSENE-D12 ** IS5 **	240	40 UG/L	1.000	4149:
1932	CI75	PERYLENE-D12 ** IS6 **	264	40 UG/L	1.000	2649:
360	CS50	2-FLUOROPHENOL ** SU1 **	112	29 %	0.667	3493
510	CS45	PHENOL-D5 ** SU2 **	99	25 %	0.944	4301:
1164	CS55	2,4,6-TRIBROMOPHENOL ** SU5 **	330	58 %	1.127	1205:
633	CS20	NITROBENZENE-D5 ** SU3 **	82	99 %	0.852	8107:
930	CS25	2-FLUOROPHENYL ** SU4 **	172	90 %	0.900	1685:
1549	CS30	TERPHENYL-D14 ** SU6 **	244	140 %	0.905	7776:
1127	CS80	DIETHYL PHTHALATE	149	1 UG/L	1.091	244

INST ID: 4530

SPECTRIX DC # -----19

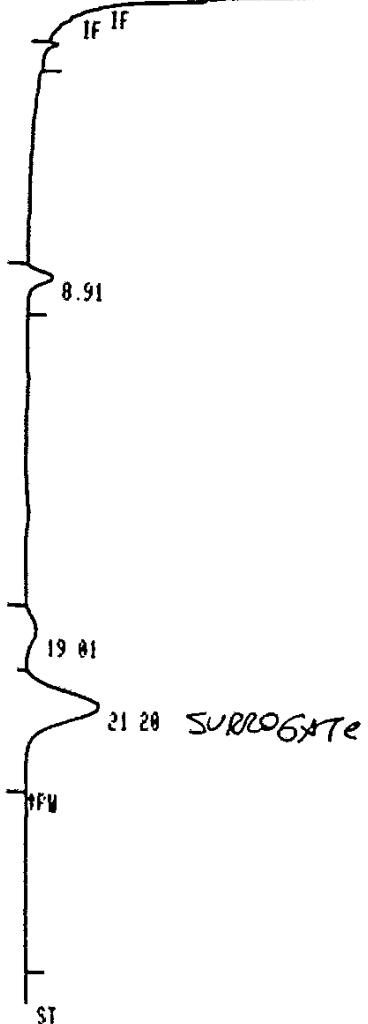
LIBRARY SEARCH
01/09/87 11:54:00 + 17:27
SAMPLE: LLW BLK
COND.: ---
ENHANCED (S 158 ZN 0T)

DATA: 4B12041C01 #1047 BASE M/Z: 73
CALI: 4B12041C01 # 3 RIC: 9439.



86-12-041-01A

R06



DC # N/A

SPECTRIX CORPORATION

CASE NO. Keystone

SAMPLE NO. LAB BLANK

PESTICIDE FRACTION

1st Column - 1D

1.8mm X 4mm ID X 6mm 00

1.5% SP 2250

1.95% SP2401

100/120 Supelcoport

Column Temp. 203 °C

Volume Inj. 1.0 ul

Instrument ID VARIAN 6c

003447

RUN # 134
WORKFILE ID: C
WORKFILE NAME:

DEC/23/86 19:45:16

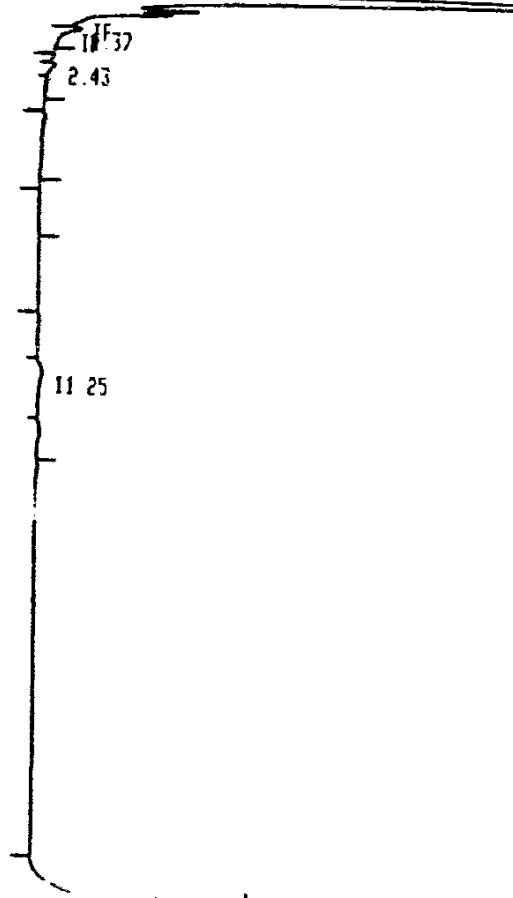
HEIGHT%

RT	HEIGHT	TYPE	AR/HT	HEIGHT%
8.91	3268	PB	0.437	23.251
19.01	1383	BY	0.886	9.364
21.20	9378	VB	0.991	66.885

TOTAL HGT= 14021
MUL FACTOR= 1.0000E+00

TAR 86-12-041-01A

N4



DC#
SPECTRIX HOUSTON

CASE NO. Keystone

SAMPLE NO. Lab. Blank.
PESTICIDE FRACTION
2ND COLUMN CONFIRMATI

2ND COLUMN II C
1.8m x2mm ID x6mm OD
3% SP-2100
100/120 SUPELCOPORT
COLUMN TEMP. 183 °C
VOLUME INJ. 1 μl
INSTRUMENT ID. 6008

M
00

JUN 4 11
WORKFILE ID: C
WORKFILE NAME:

HEIGHT%

RT	HEIGHT	TYPE	AK/HT	HEIGHT%
1.32	3185	D BB	0.007	12.890
2.43	1556	PP	0.137	5.907
11.25	1329	PY	0.648	5.045
26.83	20273	PB	1.230	76.958

TOTAL HGT= 26343
UL FACTOR= 1.0000E+00

Page 1

Received: 12/09/86

SPECTRUM 0.0

REPORT
Results by Sample

Work Order # 86-12-041

SAMPLE ID LAB BLANK (LLW)		SAMPLE # 01 FRACTIONS: A			
		Date & Time Collected not specified		Category	
AG	<10	AS	<10	BE	<3
	ug/L		ug/L		ug/L
FE	<30	HG	<0.2	NI	<29
	ug/L		ug/L		ug/L
TL	<10	ZN	<8		
	ug/L		ug/L		

GC/MS Tuning and Mass Calibration

BROMOFLUOROBENZENE

Case Number:

Inst ID: 4000
Lab ID: ES1213V01

Laboratory:
Sens Date: 12/13/86
Cali Date:

Contract:
Sens Time: 9:03:00
Analyst: CKT

Data release authorized by:

WPH

m/z	Ion Abundance Criteria	Spec # 206
50	15 to 40% of mass 95	
75	30 to 60% of mass 95	32.75
95	base peak, 100% relative abundance	43.05
96	5 to 9% of mass 95	100.00
173	less than 1% of mass 95	8.01
174	greater than 50% of mass 95	0.00
175	5 to 9% of mass 174	84.39
176	between 95% and 101% of mass 174	7.27 (8.62) 1
177	5 to 9% of mass 176	80.36 (95.22) 1 6.01 (7.48) 2

1 - value in parenthesis is % of mass 174

2 - value in parenthesis is % of mass 176

003450

Mass List
12/13/86 9:03:00 + 10:18
Sample: EFB
Conds. : ----

Data: ES1213V01 # 206 Base m/z: 95
Calib: ES1213V01 # 3 RIC: 44928.

Mass	RA	% RIC	# Inten.	Minima	Maxima	Min Inten:	0.
36. 93?	2. 93	0. 49	221.				
36. 96?F	13. 07	2. 19	985.				
37. 96?F	11. 36	1. 91	856.				
38. 99?	6. 17	1. 03	465.				
39. 93?	8. 90	1. 49	671.				
40. 99?	4. 80	0. 81	362.				
42. 00?	1. 87	0. 31	141.				
43. 02?F	7. 88	1. 32	594.				
43. 98 F	14. 52	2. 44	1094.				
45. 01	8. 01	1. 34	604.				
47. 01	4. 76	0. 80	359.				
49. 02 F	6. 40	1. 07	482.				
50. 02 F	32. 75	5. 49	2468.				
51. 04	9. 87	1. 66	744.				
55. 06	4. 01	0. 67	302.				
56. 04	3. 29	0. 55	248.				
57. 06	7. 47	1. 25	563.				
58. 06	1. 75	0. 29	132.				
61. 02	4. 46	0. 75	336.				
62. 04	4. 23	0. 71	319.				
63. 04	3. 66	0. 61	276.				
67. 10 F	1. 55	0. 26	117.				
68. 02 F	11. 45	1. 92	863.				
69. 02	11. 53	1. 93	869.				
70. 05	1. 66	0. 28	125.				
71. 07	2. 14	0. 36	161.				
73. 03	4. 60	0. 77	347.				
74. 01 F	12. 90	2. 16	972.				
75. 00 F	43. 05	7. 22	3244.				
76. 01	4. 37	0. 73	329.				
77. 02	2. 06	0. 34	155.				
78. 94	1. 73	0. 29	130.				
80. 98	2. 67	0. 45	201.				
83. 12	1. 46	0. 24	110.				
84. 99	4. 21	0. 71	317.				
87. 02 F	7. 86	1. 32	592.				
88. 02 F	7. 68	1. 29	579.				
92. 11	2. 15	0. 36	162.				
93. 08	3. 97	0. 67	299.				
94. 05 F	11. 05	1. 85	833.				
95. 02 F	100. 00	14. 77	7536.				
95. 99	8. 01	1. 34	604.				
97. 05	1. 63	0. 27	123.				
149. 03	3. 11	0. 52	234.				
173. 94 F	84. 39	14. 16	6360.				
174. 98 F	7. 27	1. 22	548.				
175. 95 F	80. 36	13. 48	6056.				
176. 98 F	6. 01	1. 01	453.				
207. 06	3. 86	0. 65	291.				

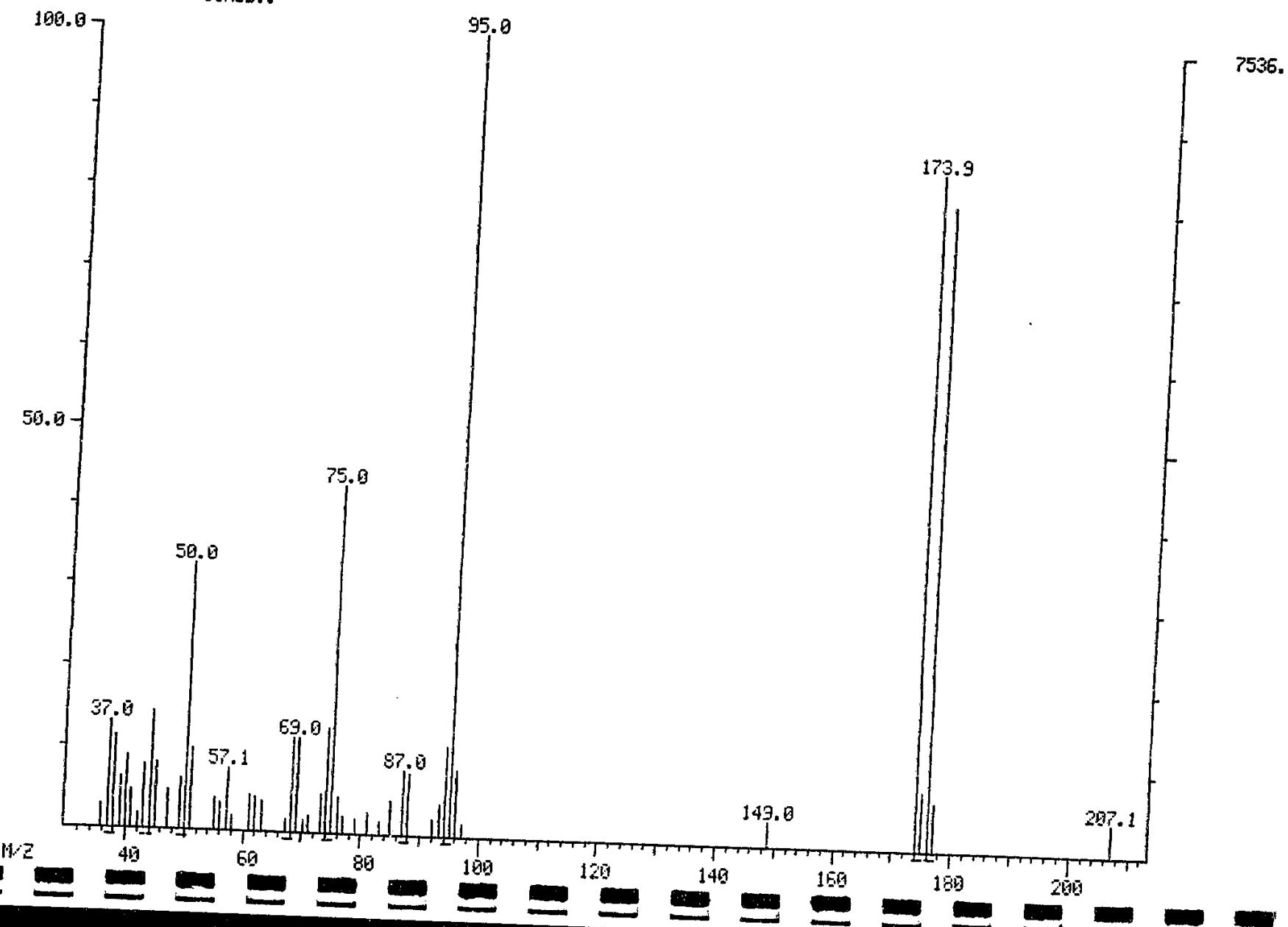
003451

003452

MASS SPECTRUM
12/13/86 9:03:00 + 10:18
SAMPLE: BFB
COND.: ----

DATA: ES1213U01 #206
CALI: ES1213U01 #3

BASE M/Z: 95
RIC: 44928.



003452

GC/MS Tuning and Mass Calibration

DECAFLUOROTRIPHENYLPHOSPHINE

Case Number:

Inst ID:

4530

Laboratory:

Lab ID:

ES0109C01

Sens Date: 01/09/87

Cali Date:

Contract:

Sens Time:

8:00:00

Analyst: GU

Data release authorized by:

GU

M/E

Ion Abundance Criteria

Spec # 395

51	30 to 60% of mass 198	47.51
68	less than 2% of mass 69	0.00 (0.00) 1
69	mass 69 relative abundance	22.28
70	less than 2% of mass 69	0.17 (0.77) 1
127	40 TO 60% of mass 198	54.92
197	less than 1% of mass 198	0.00
198	base peak, 100% relative abundance	100.00
199	5 TO 9% of mass 198	6.20
273	10 TO 30% of mass 198	21.80
365	greater than 1% of mass 198	1.50
441	less than mass 443	10.93
442	greater than 40% of mass 198	91.32
443	17 TO 23% of mass 442	16.64 (18.22) 2

1 - value in parenthesis is % of mass 69

2 - value in parenthesis is % of mass 442

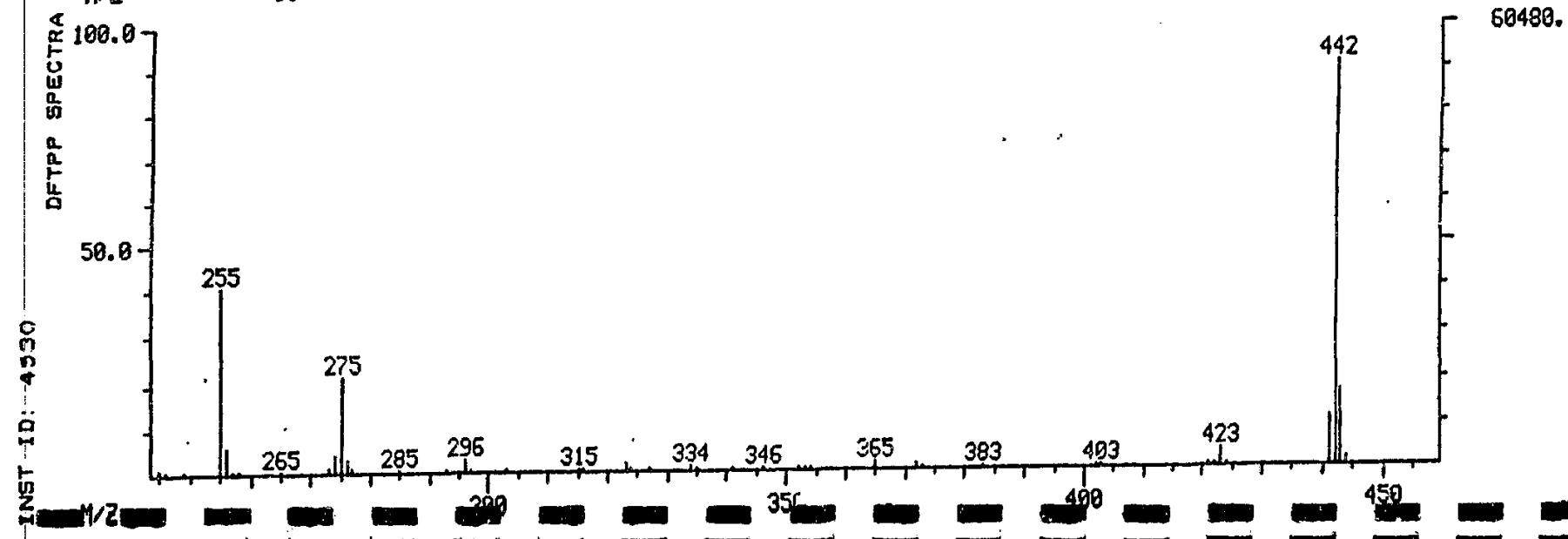
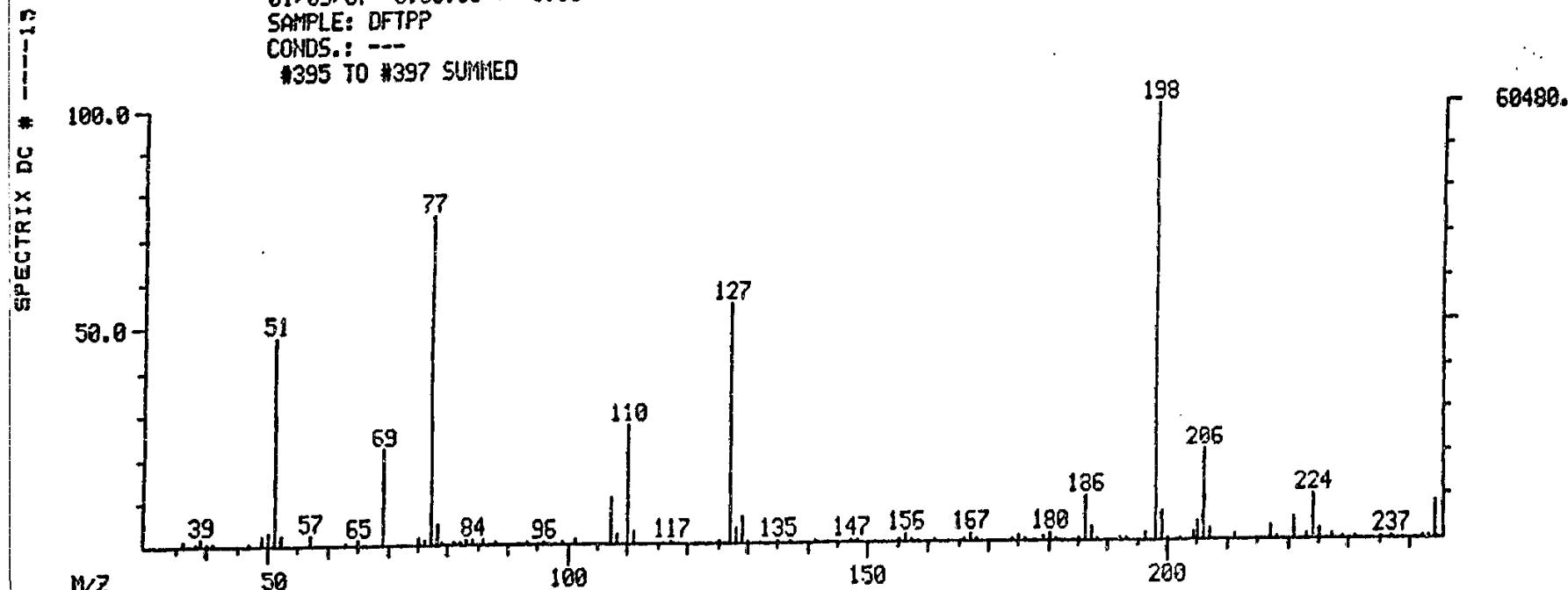
003453

003454

MASS SPECTRUM
01/09/87 8:00:00 + 6:36
SAMPLE: DFTPP
COND.: ---
#395 TO #397 SUMMED

DATA: ES0109C01 #396
CALI: ES0109C01 #3

BASE M/Z: 198
RIC: 435712.



003454

INST ID: 4530

Mass List
 01/09/87 8:00:00 + 6:36 Data: ES0109C01 # 396
 Sample: DFTPP Cali: ES0109C01 # 3
 Bonds: --- Base m/z: 198
 395 to #397 summed RIC: 435712.

SPECTRIX DC # ----15

Mass	0.00	0.	Minima Maxima	Min Inten: # 0	254.
	% RA	Inten.	Mass	% RA	Inten.
36					
445					
36?	1.09	658.	199	6.20	3752.
38?	0.45	271.	200	0.49	298.
39?	1.92	1160.	204	1.85	1120.
40	0.85	512.	205	3.97	2400.
47	0.42	257.	206	20.26	12256.
49	2.30	1394.	207	2.51	1520.
50	2.66	1608.	211	0.93	565.
51	47.51	28736.	217	2.97	1796.
52	2.27	1374.	218	0.51	308.
57	2.24	1356.	221	4.38	2652.
63	0.44	268.	223	0.91	553.
65	1.44	869.	224	10.01	6056.
69	22.28	13472.	225	2.54	1534.
75	1.70	1026.	227	1.42	856.
76	1.44	873.	229	0.72	437.
77	75.03	45376.	231	0.50	302.
78	4.80	2904.	242	0.48	292.
81	0.72	438.	243	0.54	328.
83	1.20	725.	244	8.10	4896.
84	1.28	777.	245	1.04	628.
86	0.91	551.	246	0.58	349.
101	1.17	708.	249	0.46	278.
107	10.19	6160.	255	40.58	24544.
108	2.06	1248.	256	5.70	3448.
110	27.20	16448.	257	0.44	264.
111	2.82	1706.	258	0.86	519.
117	0.43	259.	273	1.35	819.
127	54.92	33216.	274	4.03	2436.
128	3.54	2140.	275	21.80	13184.
129	6.07	3672.	276	2.75	1664.
130	0.48	289.	277	0.97	589.
135	0.63	379.	296	2.79	1686.
137	0.42	257.	297	0.45	272.
155	0.50	301.	303	0.52	313.
156	1.63	983.	323	1.88	1138.
161	0.65	393.	334	1.06	641.
167	2.02	1222.	352	0.55	333.
168	0.71	430.	354	0.75	454.
174	0.46	276.	365	1.50	905.
175	1.31	791.	372	1.01	610.
179	1.04	631.	402	0.49	296.
180	1.67	1008.	403	0.60	365.
181	0.79	479.	421	0.61	370.
185	0.66	399.	422	0.54	327.
186	9.84	5992.	423	3.96	2392.
187	2.89	1748.	424	0.74	448.
192	0.48	290.	441	10.93	6608.
193	0.66	397.	442	91.32	55232.
196	1.46	880.	443	16.64	10064.
198	100.00	60480.	444	1.57	951.

003455

GC/MS Tuning and Mass Calibration

DECAFLUOROTRIPHENYLPHOSPHINE

Case Number:

Inst ID: 4530
Lab ID: ES0111C01

Laboratory:
Sens Date: 01/11/87
Cali Date:

Contract:
Sens Time: 10:47:00
Analyst: GU

Data release authorized by:

64

M/E

Ion Abundance Criteria

Spec # 395

51	30 to 60% of mass 198	47.42
68	less than 2% of mass 69	0.00 (0.00) 1
69	mass 69 relative abundance	22.54
70	less than 2% of mass 69	0.18 (0.79) 1
127	40 TO 60% of mass 198	55.75
197	less than 1% of mass 198	0.00
198	base peak, 100% relative abundance	100.00
199	5 TO 9% of mass 198	6.12
275	10 TO 30% of mass 198	22.05
365	greater than 1% of mass 198	1.58
441	less than mass 443	11.58
442	greater than 40% of mass 198	93.15
443	17 TO 23% of mass 442	17.46 (18.74) 2

1 - value in parenthesis is % of mass 69

2 - value in parenthesis is % of mass 442

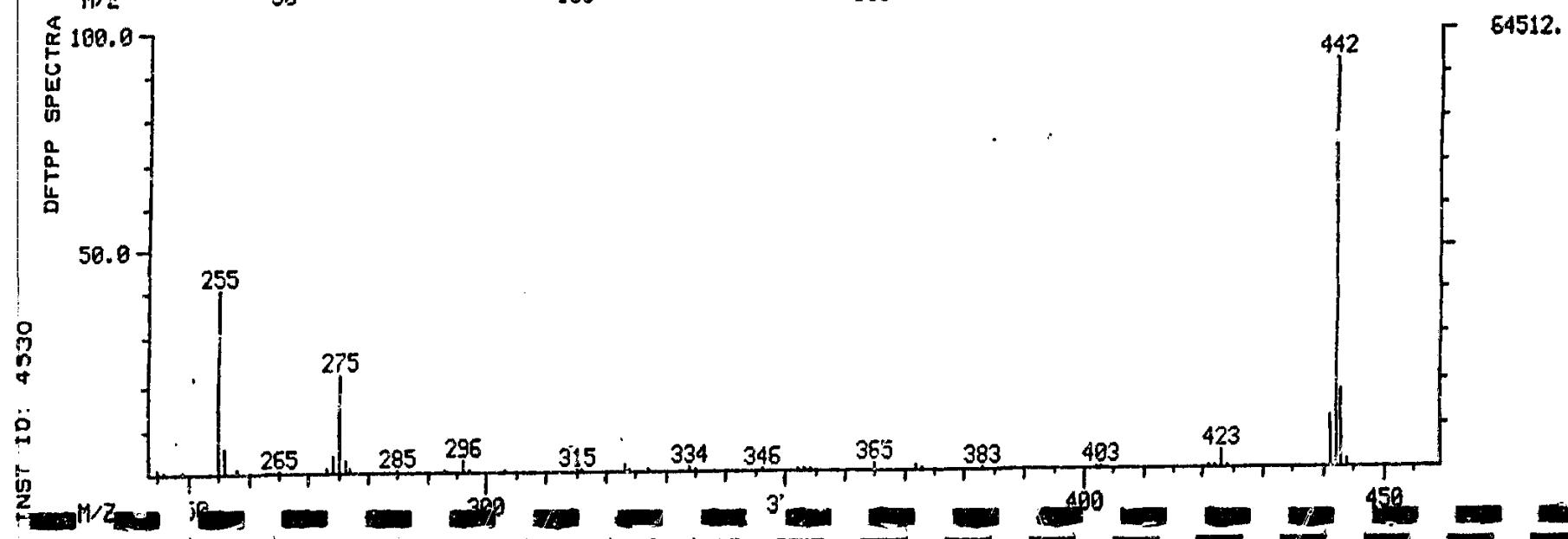
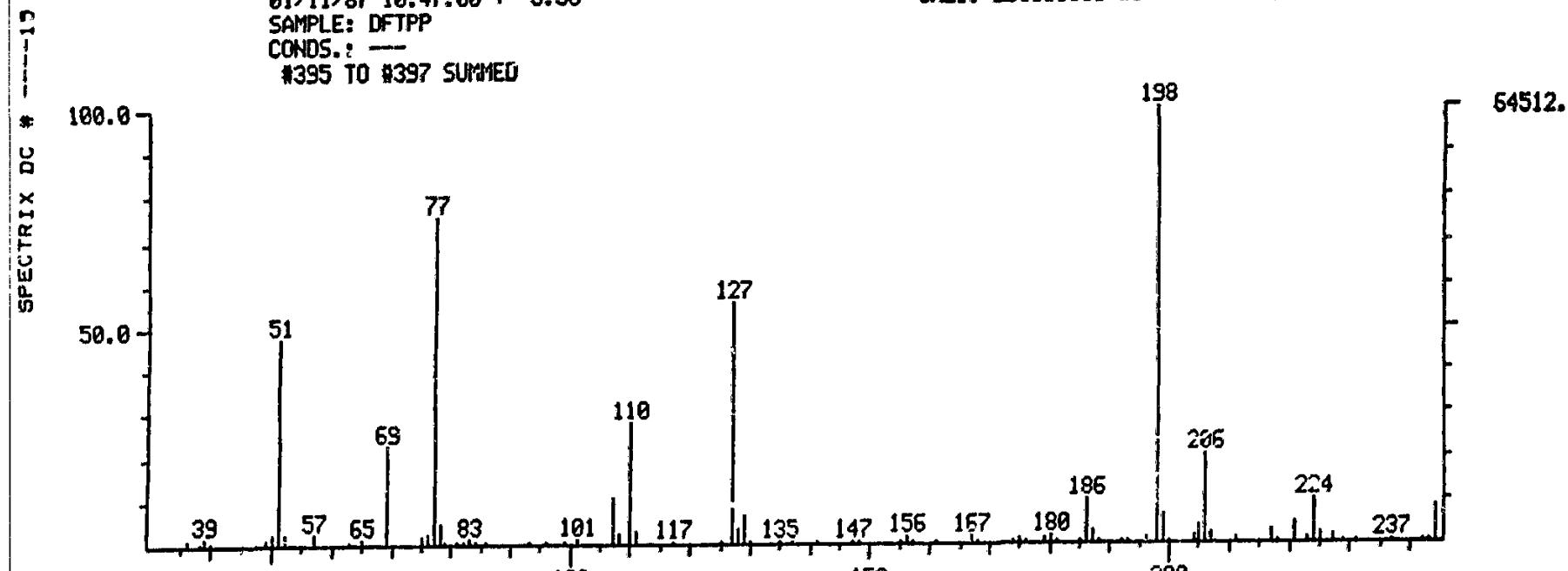
003456

003457

MASS SPECTRUM
01/11/87 10:47:00 + 6:36
SAMPLE: DFTPP
COND.: —
#395 TO #397 SUMMED

DATA: ES0111C01 #396
CALI: ES0111C01 #3

BASE M/Z: 198
RIC: 462336.



003457

INST ID: 4530

DFTPP MASS LISTING

SPECTRIX DC # ----15

Mass List
 01/11/87 10:47:00 + 6:36 Data: ES0111CO1 # 396
 Sample: DFTPP Cali: ES0111CO1 # 3
 Tnds.: ---
 395 to #397 summed

Base m/z: 198
 RIC: 462336.

Mass	0.00	0.	Minima Maxima Mass	Min Inten: # 0	243.
	% RA	Inten.		% RA	Inten.
36					
445					
36?	0.91	586.	200	0.43	279.
39?	1.93	1244.	204	1.79	1152.
40	0.68	436.	205	3.79	2448.
49	0.94	605.	206	20.26	13072.
50	2.59	1674.	207	2.46	1584.
51	47.42	30592.	211	0.90	579.
52	2.25	1432.	217	2.86	1846.
57	2.25	1454.	218	0.46	295.
63	0.42	269.	221	4.44	2864.
65	1.39	896.	223	0.91	585.
69	22.54	14544.	224	9.91	6392.
75	1.75	1130.	225	2.45	1578.
76	2.25	1454.	227	1.48	955.
77	75.30	48576.	229	0.74	475.
78	4.88	3148.	231	0.51	329.
81	0.71	461.	242	0.48	310.
83	1.16	731.	243	0.52	333.
84	0.56	360.	244	0.20	5208.
86	0.38	246.	245	1.03	665.
101	1.20	776.	246	0.54	346.
107	10.40	6712.	249	0.38	248.
108	2.10	1354.	255	40.43	26080.
110	27.68	17856.	256	5.54	3572.
111	2.86	1846.	258	0.95	411.
117	0.47	303.	273	1.31	847.
127	55.75	35948.	274	3.91	2520.
128	3.58	2308.	275	22.05	14224.
129	6.14	3964.	276	2.76	1778.
130	0.48	311.	277	0.93	603.
135	0.62	400.	293	0.49	317.
137	0.42	271.	296	2.78	1792.
141	0.40	258.	297	0.39	254.
155	0.53	339.	303	0.52	336.
156	1.63	1066.	323	1.89	1220.
161	0.64	416.	334	1.10	712.
167	2.00	1290.	352	0.56	361.
168	0.71	459.	353	0.41	262.
174	0.45	290.	354	0.74	477.
175	1.26	815.	363	1.58	1022.
179	1.10	707.	372	1.04	668.
180	1.62	1048.	402	0.46	295.
181	0.78	502.	403	0.65	422.
185	0.60	387.	421	0.62	397.
186	9.80	6320.	422	0.51	327.
187	2.77	1784.	423	4.14	2668.
192	0.47	300.	424	0.75	482.
193	0.65	420.	441	11.58	7472.
194	1.45	936.	442	93.15	60096.
198	100.00	64312.	443	17.46	11264.
199	6.12	3948.	444	1.52	978.

003500